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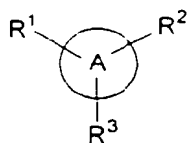
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(54) Heteroaryl-substituted pyrrole derivatives, their preparation and their therapeutic uses

(57) Compounds of formula (I):



(I)

[wherein: A is a pyrrole ring; R¹ is an optionally substituted aryl or heteroaryl group; R² is an optionally substituted nitrogen-containing heteroaryl group; R³ repre-

sents a group of the formula -X-R⁴, wherein X is a single bond or an optionally substituted alkylene, alkenylene or alkynylene group, and R⁴ is a substituted cycloalkyl group, a substituted aryl group, a substituted heterocyclyl group, an optionally substituted nitrogen-containing heterocyclyl group, a substituted heteroaryl group, an optionally substituted nitrogen-containing heteroaryl group, or -NR^aR^b, wherein each of R^a and R^b is a hydrogen atom or an alkyl, alkenyl, alkynyl, aralkyl or alkyl-sulfonyl group; PROVIDED THAT said substituents R¹ and R³ are bonded to the two atoms of said pyrrole ring which are adjacent to the atom of the pyrrole ring to which said substituent R² is bonded] have excellent inhibitory activity against the production of inflammatory cytokines.

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Description

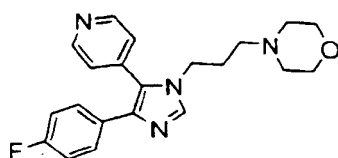
[0001] The present invention relates to a series of heteroaryl-substituted pyrrole derivatives which have excellent inhibitory activity against the production of inflammatory cytokines such as interleukin (IL)-1, IL-6 and IL-8 and tumor necrosis factor (TNF), particularly IL-1 and TNF. As a consequence, the compounds of the present invention have valuable anti-pyretic, analgesic, anti-viral and anti-inflammatory activity and are useful in the prophylaxis and treatment of autoimmune diseases such as chronic rheumatism, bone diseases such as osteoporosis and the many other diseases in which the above-described inflammatory cytokines take part. The invention also provides compositions containing these novel compounds, uses of said novel compounds and processes for their preparation.

[0002] Non-steroidal anti-inflammatory drugs (NSAIDs) have been widely used for the treatment and prophylaxis of various inflammatory diseases and in pain relief because they have, as their main pharmacological activity, anti-pyretic, analgesic, and anti-inflammatory activity which is based on their ability to inhibit the biosynthesis of prostaglandin (PG) through the inhibition of cyclooxygenase activity. Another class of compounds commonly used for the treatment of rheumatoid arthritis is the disease-modifying anti-rheumatic drugs (DMARDs), examples of which include methotrexate and sulphasalazine. This is a wide class of drugs in which the compounds have no common mechanism of action. For the treatment of chronic rheumatism, NSAIDs are used nosotopically and DMARDs are used etiotropically. There are a number of problems associated with these classes of drugs. Conventional NSAIDs can induce undesirable side effects including gastrointestinal disorders such as gastric ulcers and renal disorders, resulting in difficulties for any patient who has to take such a drug for an extended period of time. DMARDs can also induce undesirable side effects including nausea and diarrhoea and, furthermore, they have not yet been clearly shown to exhibit a stable, long-lasting effect.

[0003] A class of active substances generally called cytokines, which are produced in the body by immunocytes, has recently been found. One group of cytokines is known as the inflammatory cytokines and it includes interleukin (IL)-1, IL-6 and IL-8 and tumor necrosis factor (TNF). The inflammatory cytokines have been demonstrated to play a major role in a number of biological processes. These include action as an inflammatory mediator through the stimulation of the arachidonic acid metabolic pathway leading to the production of PG, the migration of leukocytes, the production of acute phase protein, and activation of osteoclasts.

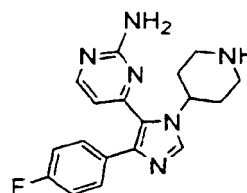
[0004] It is believed that the inflammatory cytokines are associated with many diseases including inflammatory diseases and the induction of bone resorption. Due to their mechanism of action, which is different from that of conventional drugs such as those described above, compounds which are able to inhibit the production of inflammatory cytokines are expected to provide an improved new generation of antipyretic, analgesic and anti-inflammatory drugs and medicaments for the treatment of autoimmune diseases such as chronic rheumatism, bone diseases such as osteoporosis and the many other diseases in which the above-described inflammatory cytokines are believed to take part.

[0005] Compounds which are said to demonstrate inhibitory activity against the production of inflammatory cytokines include various heteroaryl compounds [see, for example, WO 96/21452, WO 97/5877, WO 97/23479 and J. Med. Chem., 39, 3929-3937 (1996)]. Examples of compounds of this type include the following:

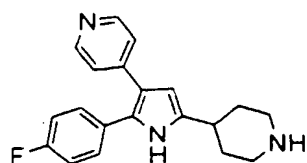


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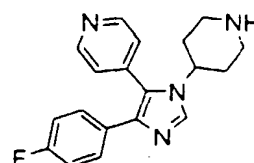
J. Med. Chem., 39,
3929-3937 (1996)



WO97/23479
Compound of
Example 6



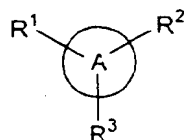
WO97/5877
Compound of
Example 4



WO96/21452
Compound of
Example 23

[0006] There is, however, a need for further compounds having improved activity, pharmacokinetics and safety.

[0007] The compounds of the present invention are compounds of the following formula (I), and pharmacologically acceptable salts, esters or other derivatives thereof:



(I)

wherein:

A represents a pyrrole ring;

R¹ represents an aryl group defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below, or a heteroaryl group defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below;

R² represents a heteroaryl group defined below having at least one ring nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below;

R³ represents a group of the formula -X-R⁴, wherein:

X is selected from the group consisting of single bonds,

lower alkylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below,

lower alkenylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below, and

lower alkynylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below; and

R⁴ is selected from the group consisting of

cycloalkyl groups defined below which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below and which may option-

ally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below,

aryl groups defined below which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and

Substituent group γ defined below and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below,

heterocyclyl groups defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below and which are substituted with at least one substituent selected from the group consisting of Substituent group β

defined below and Substituent group γ defined below,

heterocyclyl groups defined below having at least one nitrogen atom, said heterocyclyl groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below,

heteroaryl groups defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below,

heteroaryl groups defined below having at least one nitrogen atom, said heteroaryl groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below and

Substituent group δ defined below; and groups of formula $-NR^aR^b$, wherein R^a and R^b are the same or different from each other and each is independently selected from the group consisting of hydrogen atoms, lower alkyl groups defined below, lower alkenyl groups defined below, lower alkynyl groups defined below, aralkyl groups defined below and lower alkylsulfonyl groups defined below;

PROVIDED THAT said substituents R^1 and R^3 are bonded to the two atoms of said pyrrole ring which are adjacent to the atom of the pyrrole ring to which said substituent R^2 is bonded;

Substituent group α comprises hydroxyl groups, nitro groups, cyano groups, halogen atoms, lower alkoxy groups defined below, halogeno lower alkoxy groups defined below, lower alkylthio groups defined below, and halogeno lower alkylthio groups defined below;

Substituent group β comprises groups of formula $-NR^cR^d$, wherein R^c and R^d are the same or different from each other and each is independently selected from the group consisting of hydrogen atoms, lower alkyl groups defined below, lower alkenyl groups defined below, lower alkynyl groups defined below, aralkyl groups defined below and lower alkylsulfonyl groups, or

R^c and R^d , together with the nitrogen atom to which R^c and R^d are bonded, form a heterocyclyl group defined below;

Substituent group γ comprises lower alkyl groups defined below which are substituted with a group of formula $-NR^cR^d$, wherein R^c and R^d are as defined above; and

Substituent group δ comprises lower alkyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, lower alkenyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, lower alkynyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, aralkyl groups defined below and cycloalkyl groups defined below.

[0008] The present invention also provides a pharmaceutical composition comprising an effective amount of a pharmacologically active compound together with a carrier or diluent therefor, wherein said pharmacologically active compound is a compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof.

[0009] The present invention also provides a compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof for use as a medicament.

[0010] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for inhibiting the production of inflammatory cytokines in a mammal, which may be human.

[0011] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for inhibiting bone resorption in a mammal, which may be human.

[0012] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of inflammatory diseases in a mammal, which may be human.

[0013] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of viral diseases in a mammal, which may be human.

[0014] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for relieving pain or pyrexia in a mammal, which may be human.

[0015] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of chronic rheumatoid arthritis in a mammal, which may be human.

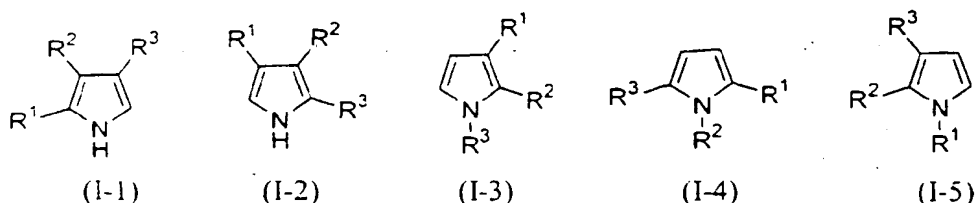
[0016] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of osteoarthritis in a mammal, which may be human.

[0017] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of cancer in a mammal, which may be human.

[0018] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of hepatitis in a mammal, which may be human.

[0019] The present invention also provides the use of at least one compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of a disease selected from the group consisting of allergic diseases, septicæmia, psoriasis, osteoporosis, ulcerative colitis, diabetes, nephritis, ischemic heart disease, Alzheimer's disease and arteriosclerosis in a mammal, which may be human.

[0020] The requirement that the substituents R^1 and R^3 must be bonded to the two atoms of the pyrrole ring which are adjacent to the atom of the pyrrole ring to which the substituent R^2 is bonded in the compounds of the above formula (I) means that the compounds of formula (I) are selected from compounds of the following formulae (I-1) to (I-5):



wherein R^1 , R^2 and R^3 are as defined above.

[0021] Where R^1 represents an aryl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above or where R^4 represents an aryl group which is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, said aryl groups are aromatic hydrocarbon groups having from 6 to 14 carbon atoms in one or more rings, preferably from 6 to 10 carbon atoms, and examples include phenyl, naphthyl, phenanthryl and anthracenyl groups. Of these, we prefer phenyl and naphthyl groups, most preferably phenyl groups.

[0022] The aryl groups defined and exemplified above may be fused with a cycloalkyl group having from 3 to 10 carbon atoms. Examples of such a fused ring group include 5-indanyl groups.

[0023] Where R^1 represents an aryl group which is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above, it is preferably an aryl group substituted with 1 to 4 substituents selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , and more preferably it is an aryl group substituted with 1 to 3 substituents selected from Substituent group α , Substituent group β , Substituent group γ and Substituent group δ . Examples of such substituted aryl groups include 4-fluorophenyl,

3-fluorophenyl, 4-chlorophenyl, 3-chlorophenyl, 3,4-difluorophenyl, 3,4-dichlorophenyl, 3,4,5-trifluorophenyl, 3-chloro-4-fluorophenyl, 3-difluoromethoxyphenyl, 3-trifluoromethoxyphenyl and 3-trifluoromethylphenyl groups.

[0024] Where R⁴ represents an aryl group which is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, it is preferably an aryl group substituted with a group selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with one or two groups selected from the group consisting of Substituent group α .

[0025] Examples of such substituted aryl groups include 2-aminophenyl, 3-aminophenyl, 4-aminophenyl, 2-aminomethylphenyl, 3-aminomethylphenyl, 4-aminomethylphenyl, 2-(2-aminoethyl)phenyl, 3-(2-aminoethyl)phenyl, 4-(2-aminoethyl)phenyl, 2-methylaminophenyl, 3-methylaminophenyl, 4-methylaminophenyl, 2-(dimethylamino)phenyl, 3-(dimethylamino)phenyl, 4-(dimethylamino)phenyl, 2-methylaminomethylphenyl, 3-methylaminomethylphenyl, 4-methylaminomethylphenyl, 2-(dimethylaminomethyl)phenyl, 3-(dimethylaminomethyl)phenyl, 4-(dimethylaminomethyl)phenyl, 3-amino-4-fluorophenyl, 3-amino-5-fluorophenyl, 2-aminomethyl-4-fluorophenyl, 3-amino-5-chlorophenyl, 2-aminomethyl-3-chlorophenyl, 3-amino-5-difluoromethoxyphenyl and 2-aminomethyl-3-trifluoromethoxyphenyl groups.

[0026] Where R¹ represents a heteroaryl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above, or where R⁴ represents a heteroaryl group which is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, said heteroaryl groups are 5- to 7-membered aromatic heterocyclic groups containing from 1 to 3 heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms. Examples of such heteroaryl groups include furyl, thienyl, pyrrolyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, triazolyl, tetrazolyl, thiadiazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups. We prefer 5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, examples of which include furyl, thienyl, pyrrolyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups.

[0027] Where R¹ represents a heteroaryl group, 5- or 6-membered aromatic heterocyclic groups containing 1 or 2 heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms are more preferred, and furyl, thienyl, pyridyl and pyrimidinyl groups are particularly preferred.

[0028] Where R⁴ represents a heteroaryl group, 5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms and optionally containing a further heteroatom selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms are preferred, examples of which include pyrrolyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups. Of these, 5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, such as imidazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups are preferred, and pyridyl and pyrimidinyl groups are particularly preferred.

[0029] The heteroaryl groups defined and exemplified above may be fused with another cyclic group selected from the group consisting of aryl groups defined above and cycloalkyl groups having from 3 to 10 carbon atoms. Examples of such a fused heteroaryl group include indolyl, benzofuranyl, benzothienyl, quinolyl, isoquinolyl, quinazolyl, tetrahydroquinolyl and tetrahydroisoquinolyl groups.

[0030] Where R¹ represents a heteroaryl group which is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , said heteroaryl group is preferably a heteroaryl group substituted with from 1 to 3 substituents selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , and more preferably it is a heteroaryl group substituted with one or two substituents selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ . Examples of such substituted heteroaryl groups include 5-fluoro-2-furyl, 4-chloro-2-thienyl, 5-difluoromethoxy-3-furyl, 5-trifluoromethyl-3-thienyl and 5-fluoro-2-oxazolyl groups.

[0031] Where R⁴ represents a heteroaryl group which is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, it is preferably a heteroaryl group substituted with a substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with a substituent selected from Substituent group α .

[0032] Examples of such substituted heteroaryl groups include 5-amino-2-furyl, 5-aminomethyl-2-furyl, 5-methylaminomethyl-2-furyl, 5-dimethylaminomethyl-2-furyl, 5-amino-2-thienyl, 5-aminomethyl-2-thienyl, 5-methylaminomethyl-

2-thienyl, 5-dimethylaminomethyl-2-thienyl, 5-amino-2-oxazolyl, 5-aminomethyl-2-oxazolyl, 2-amino-4-pyridyl, 2-amino-4-pyrimidinyl, 2-methylamino-4-pyridyl, 2-methylamino-4-pyrimidinyl, 2-benzylamino-4-pyridyl, 2-benzylamino-4-pyrimidinyl, 2-(α -methylbenzylamino)-4-pyridyl, 2-(α -methylbenzylamino)-4-pyrimidinyl, 5-amino-4-fluoro-2-furyl, 5-aminomethyl-4-fluoro-2-furyl, 5-amino-4-fluoro-2-thienyl, 5-aminomethyl-4-fluoro-2-thienyl, 4-amino-5-difluoromethoxy-2-furyl, 4-aminomethyl-5-difluoromethoxy-2-furyl, 4-amino-5-difluoromethoxy-2-thienyl and 4-aminomethyl-5-difluoromethoxy-2-thienyl groups.

[0033] Where R^2 represents a heteroaryl group having at least one ring nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above or R^4 represents a heteroaryl group having at least one ring nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, said heteroaryl groups are 5- to 7-membered aromatic heterocyclic groups containing one or two nitrogen atoms and optionally containing one or two further heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms. Examples of such groups include pyrrolyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, triazolyl, tetrazolyl, thiadiazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups. Of these, we prefer 5- or 6-membered aromatic heterocyclic groups containing one nitrogen atom and optionally containing one further heteroatom selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, examples of which include pyrrolyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups. 5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, such as imidazolyl, pyridyl, pyridazinyl, pyrimidinyl and pyrazinyl groups are more preferred and pyridyl and pyrimidinyl groups are particularly preferred.

[0034] Where R^2 represents a heteroaryl group having at least one ring nitrogen atom, 4-pyridyl and 4-pyrimidinyl groups are most preferred.

[0035] Where R^2 represents a heteroaryl group having at least one ring nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , said heteroaryl group is preferably a group substituted with 1 to 3 substituents selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , more preferably it is a heteroaryl group substituted with one or two substituents selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , still more preferably it is a heteroaryl group substituted with one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , and particularly preferably it is a 4-pyridyl or 4-pyrimidinyl group which is substituted with one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ at the 2-position of said group. Most preferably, said heteroaryl group is a 4-pyridyl or 4-pyrimidinyl group which is substituted at the 2-position with one substituent selected from the group consisting of Substituent group β and Substituent group γ . Examples of such substituted heteroaryl groups include 2-amino-4-pyridyl, 2-amino-4-pyrimidinyl, 2-methylamino-4-pyridyl, 2-methylamino-4-pyrimidinyl, 2-methoxy-4-pyridyl, 2-methoxy-4-pyrimidinyl, 2-benzylamino-4-pyridyl, 2-benzylamino-4-pyrimidinyl, 2-(α -methylbenzylamino)-4-pyridyl and 2-(α -methylbenzylamino)-4-pyrimidinyl groups.

[0036] Where R^4 represents a heteroaryl group having at least one nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, it is preferably a heteroaryl group substituted with 1 or 2 substituents selected from the group consisting of Substituent group α and Substituent group δ , and more preferably it is a heteroaryl group substituted with a substituent selected from the group consisting of Substituent group α and Substituent group δ .

[0037] Examples of such substituted heteroaryl groups include 2-methyl-4-pyridyl, 2-ethyl-4-pyridyl, 2-benzyl-4-pyridyl, 2-phenethyl-4-pyridyl, 2-fluoro-4-pyridyl, 2,6-difluoro-4-pyridyl and 2,3,5,6-tetrafluoro-4-pyridyl groups.

[0038] Where X represents a lower alkylene group which may optionally be substituted with at least one substituent selected from Substituent group α defined above, said lower alkylene group is a straight or branched alkylene group having from 1 to 6 carbon atoms, examples of which include methylene, ethylene, trimethylene, propylene, tetramethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethylethylene, pentamethylene, 1,1-dimethyltrimethylene and hexamethylene groups. Of these, straight or branched alkylene groups having from 1 to 4 carbon atoms are preferred, straight or branched alkylene groups having from 1 to 3 carbon atoms are more preferred, and methylene, ethylene and trimethylene groups are most preferred.

[0039] Where X represents a lower alkenylene group which may optionally be substituted with at least one substituent selected from Substituent group α defined above, said lower alkenylene group is a straight or branched alkenylene group having from 2 to 6 carbon atoms, examples of which include vinylene, 1-methylvinylene, propenylene, 1-butenylene, 2-butenylene, 1-pentenylene and 2-pentenylene groups. Of these, straight or branched alkenylene groups having from 2 to 4 carbon atoms are preferred, vinylene, propenylene and butenylene groups are more preferred, and vinylene and propenylene groups are most preferred.

[0040] Where X represents a lower alkynylene group which may optionally be substituted with at least one substituent selected from Substituent group α defined above, said lower alkynylene group is a straight or branched alkynylene group having from 2 to 6 carbon atoms. Of these, straight or branched alkynylene groups having from 2 to 4 carbon atoms are preferred. ethynylene, propynylene, 1-butylnylene and 2-butylnylene groups are more preferred, and ethynylene and propynylene groups are most preferred.

[0041] Where R^4 represents a cycloalkyl group which is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above or where the substituent from Substituent group δ represents a cycloalkyl group, said cycloalkyl group has from 3 to 7 carbon atoms, examples of which include cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl and cycloheptyl groups. Of these, cycloalkyl groups having from 3 to 6 carbon atoms are preferred.

[0042] Where R^4 represents a heterocyclyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and is substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above, said heterocyclyl group is a non-aromatic heterocyclic group having from 4 to 14 ring atoms in one or more rings, at least one of said ring atoms being a heteroatom selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms. It is preferably a 4- to 12-membered non-aromatic heterocyclic group (and more preferably a 4- to 10-membered non-aromatic heterocyclic group) containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, and more preferably it is a 4- to 12-membered non-aromatic heterocyclic group (preferably a 4- to 10-membered non-aromatic heterocyclic group) containing one nitrogen atom and optionally containing one further heteroatom selected from the group consisting of oxygen atoms, sulfur atoms and nitrogen atoms. Examples of such a group include azetidiny, pyrrolidiny, pyrroliny, imidazolidiny, imidazoliny, pyrazolidiny, pyrazoliny, oxazolidiny, thiazolidiny, piperidyl, tetrahydropyridyl, dihydropyridyl, piperaziny, morpholiny, thiomorpholiny, homopiperidyl, quinuclidiny, quinuclidiny, octahydroindoliziny, hexahydroindoliziny, octahydroquinoliziny, hexahydroquinoliziny, 8-azabicyclo[3.2.1]octany, 8-azabicyclo[3.2.1]octeny, 9-azabicyclo[3.3.1]nonany and 9-azabicyclo[3.3.1]noneny groups, of which piperidyl, tetrahydropyridyl, homopiperidyl, quinuclidiny, quinuclidiny, octahydroindoliziny, hexahydroindoliziny, octahydroquinoliziny, hexahydroquinoliziny, azabicyclo[3.2.1]octany, azabicyclo[3.2.1]octeny, azabicyclo[3.3.1]nonany and azabicyclo[3.3.1]noneny groups are preferred, tetrahydropyridyl, quinuclidiny, hexahydroindoliziny, hexahydroquinoliziny, azabicyclo[3.2.1]octeny and azabicyclo[3.3.1]noneny groups are more preferred, and tetrahydropyridyl, hexahydroindoliziny and hexahydroquinoliziny groups are most preferred.

[0043] The heterocyclyl groups defined and exemplified above may be fused with another cyclic group selected from the group consisting of aryl groups defined above and heteroaryl groups defined above. Examples of such a fused heterocyclyl group include tetrahydroquinoliny, tetrahydroisoquinoliny, chromany, indoliny and isoindoliny groups.

[0044] Where R^4 represents a heterocyclyl group having at least one ring nitrogen atom in which said heterocyclyl group may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, said heterocyclyl group is a 4- to 12-membered non-aromatic heterocyclic group containing one nitrogen atom and optionally containing one further heteroatom selected from the group consisting of oxygen atoms, sulfur atoms and nitrogen atoms. Preferably, it is a 4- to 10-membered non-aromatic heterocyclic group containing one nitrogen atom and optionally containing one further heteroatom selected from the group consisting of oxygen atoms, sulfur atoms and nitrogen atoms. Examples of such a group include azetidiny, pyrrolidiny, pyrroliny, imidazolidiny, imidazoliny, pyrazolidiny, pyrazoliny, oxazolidiny, thiazolidiny, piperidyl, tetrahydropyridyl, dihydropyridyl, piperaziny, morpholiny, thiomorpholiny, homopiperidyl, quinuclidiny, quinuclidiny, octahydroindoliziny, hexahydroindoliziny, octahydroquinoliziny, hexahydroquinoliziny, 8-azabicyclo[3.2.1]octany, 8-azabicyclo[3.2.1]octeny, 9-azabicyclo[3.3.1]nonany and 9-azabicyclo[3.3.1]noneny groups, of which piperidyl, tetrahydropyridyl, homopiperidyl, quinuclidiny, quinuclidiny, octahydroindoliziny, hexahydroindoliziny, octahydroquinoliziny, hexahydroquinoliziny, azabicyclo[3.2.1]octany, azabicyclo[3.2.1]octeny, azabicyclo[3.3.1]nonany and azabicyclo[3.3.1]noneny groups are preferred, tetrahydropyridyl, quinuclidiny, hexahydroindoliziny, hexahydroquinoliziny, azabicyclo[3.2.1]octeny and azabicyclo[3.3.1]noneny groups are more preferred, and tetrahydropyridyl, hexahydroindoliziny and hexahydroquinoliziny groups are most preferred.

[0045] The heterocyclyl groups having at least one ring nitrogen atom defined and exemplified above may be fused with another cyclic group selected from the group consisting of aryl groups defined above and heteroaryl groups defined above. Examples of such fused heterocyclyl groups include tetrahydroquinoliny, tetrahydroisoquinoliny, indoliny and isoindoliny groups.

[0046] The lower alkyl groups in the definition of substituents R^a , R^b , R^c and R^d , the lower alkyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ and the lower alkyl moiety of the lower alkyl groups substituted with a group of formula $-NR^cR^d$ in the definition

of Substituent group γ are straight or branched alkyl groups having from 1 to 6 carbon atoms. Examples of said lower alkyl groups include methyl, ethyl, propyl, isopropyl, n-butyl, isobutyl, s-butyl, tert-butyl, n-pentyl, isopentyl, 2-methylbutyl, neopentyl, 1-ethylpropyl, n-hexyl, isohexyl, 4-methylpentyl, 3-methylpentyl, 2-methylpentyl, 1-methylpentyl, 3,3-dimethylbutyl, 2,2-dimethylbutyl, 1,1-dimethylbutyl, 1,2-dimethylbutyl, 1,3-dimethylbutyl, 2,3-dimethylbutyl and 2-ethylbutyl groups. Alkyl groups having from 1 to 4 carbon atoms are preferred, methyl, ethyl and propyl groups are more preferred, and methyl and ethyl groups are most preferred.

[0047] The lower alkenyl groups in the definition of substituents R^a , R^b , R^c and R^d and the lower alkenyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ are straight or branched alkenyl groups having from 2 to 6 carbon atoms. Examples of said lower alkenyl groups include vinyl, 2-propenyl, 1-methyl-2-propenyl, 2-methyl-2-propenyl, 2-ethyl-2-propenyl, 2-butenyl, 1-methyl-2-butenyl, 2-methyl-2-butenyl, 1-ethyl-2-butenyl, 3-butenyl, 1-methyl-3-butenyl, 2-methyl-3-butenyl, 1-ethyl-3-butenyl, 2-pentenyl, 1-methyl-2-pentenyl, 2-methyl-2-pentenyl, 3-pentenyl, 1-methyl-3-pentenyl, 2-methyl-3-pentenyl, 4-pentenyl, 1-methyl-4-pentenyl, 2-methyl-4-pentenyl, 2-hexenyl, 3-hexenyl, 4-hexenyl and 5-hexenyl groups. Alkenyl groups having from 2 to 4 carbon atoms are preferred, and alkenyl groups having 2 or 3 carbon atoms are most preferred.

[0048] The lower alkynyl groups in the definition of substituents R^a , R^b , R^c and R^d and the lower alkynyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ are straight or branched alkynyl groups having from 2 to 6 carbon atoms. Examples of said lower alkynyl groups include ethynyl, 2-propynyl, 1-methyl-2-propynyl, 2-butylnyl, 1-methyl-2-butylnyl, 1-ethyl-2-butylnyl, 3-butylnyl, 1-methyl-3-butylnyl, 2-methyl-3-butylnyl, 1-ethyl-3-butylnyl, 2-pentylnyl, 1-methyl-2-pentylnyl, 3-pentylnyl, 1-methyl-3-pentylnyl, 2-methyl-3-pentylnyl, 4-pentylnyl, 1-methyl-4-pentylnyl, 2-methyl-4-pentylnyl, 2-hexynyl, 3-hexynyl, 4-hexynyl and 5-hexynyl groups. Alkynyl groups having from 2 to 4 carbon atoms are preferred, and alkynyl groups having 2 or 3 carbon atoms are most preferred.

[0049] The aralkyl group in the definition of R^a , R^b , R^c , R^d and Substituent group δ is a lower alkyl group as defined above which is substituted with at least one aryl group as defined above. Examples of said aralkyl group include benzyl, phenanthrenylmethyl, anthracenylmethyl, α -naphthylmethyl, β -naphthylmethyl, diphenylmethyl, triphenylmethyl, α -naphthyldiphenylmethyl, 9-anthrylmethyl, piperonyl, 1-phenethyl, 2-phenethyl, 1-naphthylethyl, 2-naphthylethyl, 1-phenylpropyl, 2-phenylpropyl, 3-phenylpropyl, 1-naphthylpropyl, 2-naphthylpropyl, 3-naphthylpropyl, 1-phenylbutyl, 2-phenylbutyl, 3-phenylbutyl, 4-phenylbutyl, 1-naphthylbutyl, 2-naphthylbutyl, 3-naphthylbutyl, 4-naphthylbutyl, 1-phenylpentyl, 2-phenylpentyl, 3-phenylpentyl, 4-phenylpentyl, 5-phenylpentyl, 1-naphthylpentyl, 2-naphthylpentyl, 3-naphthylpentyl, 4-naphthylpentyl, 5-naphthylpentyl, 1-phenylhexyl, 2-phenylhexyl, 3-phenylhexyl, 4-phenylhexyl, 5-phenylhexyl, 6-phenylhexyl, 1-naphthylhexyl, 2-naphthylhexyl, 3-naphthylhexyl, 4-naphthylhexyl, 5-naphthylhexyl and 6-naphthylhexyl groups, of which benzyl, phenanthrenylmethyl, anthracenylmethyl, α -naphthylmethyl, β -naphthylmethyl, diphenylmethyl, triphenylmethyl, 9-anthrylmethyl, piperonyl, 1-phenethyl, 2-phenethyl, 1-phenylpropyl, 2-phenylpropyl, 3-phenylpropyl, 1-phenylbutyl, 2-phenylbutyl, 3-phenylbutyl and 4-phenylbutyl are preferred.

[0050] The aryl moiety of the aralkyl groups defined and exemplified above may be substituted with from 1 to 3 groups selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above. Examples of said substituted aralkyl groups include aralkyl groups substituted with at least one halogen atom, examples of which include 2-fluorobenzyl, 3-fluorobenzyl, 4-fluorobenzyl, 2-chlorobenzyl, 3-chlorobenzyl, 4-chlorobenzyl, 2-bromobenzyl, 3-bromobenzyl, 4-bromobenzyl, 3,5-difluorobenzyl, 2,5-difluorophenethyl, 2,6-difluorobenzyl, 2,4-difluorophenethyl, 3,5-dibromobenzyl, 2,5-dibromophenethyl, 2,6-dichlorobenzyl, 2,4-dichlorophenethyl, 2,3,6-trifluorobenzyl, 2,3,4-trifluorophenethyl, 3,4,5-trifluorobenzyl, 2,5,6-trifluorophenethyl, 2,4,6-trifluorobenzyl, 2,3,6-tribromophenethyl, 2,3,4-tribromobenzyl, 3,4,5-tribromophenethyl, 2,5,6-trichlorobenzyl, 2,4,6-trichlorophenethyl, 1-fluoro-2-naphthylmethyl, 2-fluoro-1-naphthylethyl, 3-fluoro-1-naphthylmethyl, 1-chloro-2-naphthylethyl, 2-chloro-1-naphthylmethyl, 3-bromo-1-naphthylethyl, 3,8-difluoro-1-naphthylmethyl, 2,3-difluoro-1-naphthylethyl, 4,8-difluoro-1-naphthylmethyl, 5,6-difluoro-1-naphthylethyl, 3,3-dichloro-1-naphthylmethyl, 2,3-dichloro-1-naphthylethyl, 4,8-dibromo-1-naphthylmethyl, 5,6-dibromo-1-naphthylethyl, 2,3,6-trifluoro-1-naphthylmethyl, 2,3,4-trifluoro-1-naphthylethyl, 3,4,5-trifluoro-1-naphthylmethyl, 4,5,6-trifluoro-1-naphthylethyl, 2,4,8-trifluoro-1-naphthylmethyl, bis(2-fluorophenyl)methyl, 3-fluorophenylphenylmethyl, bis(4-fluorophenyl)methyl, 4-fluorophenylphenylmethyl, bis(2-chlorophenyl)methyl, bis(3-chlorophenyl)methyl, bis(4-chlorophenyl)methyl, 4-chlorophenylphenylmethyl, 2-bromophenylphenylmethyl, 3-bromophenylphenylmethyl, bis(4-bromophenyl)methyl, bis(3,5-difluorophenyl)methyl, bis(2,5-difluorophenyl)methyl, bis(2,6-difluorophenyl)methyl, 2,4-difluorophenylphenylmethyl, bis(3,5-dibromophenyl)methyl, 2,5-dibromophenylphenylmethyl, 2,6-dichlorophenylphenylmethyl, bis(2,4-dichlorophenyl)methyl and bis(2,3,6-trifluorophenyl)methyl groups; aralkyl groups substituted with at least one halogeno lower alkyl group (as defined below), examples of which include 2-trifluoromethylbenzyl, 3-trifluoromethylphenethyl, 4-trifluoromethylbenzyl, 2-trichloromethylphenethyl, 3-dichloromethylbenzyl, 4-trichloromethylphenethyl, 2-tribromomethylbenzyl, 3-dibromomethylphenethyl, 4-dibromomethylbenzyl, 3,5-bistrifluoromethylphenethyl, 2,5-bistrifluoromethylbenzyl, 2,6-bistrifluoromethylphenethyl, 2,4-bistrifluoromethylbenzyl, 3,5-bistribromomethylphenethyl,

2,5-bisdibromomethylbenzyl, 2,6-bisdichloromethylphenethyl, 2,4-bisdichloromethylbenzyl, 2,3,6-tris(trifluoromethyl)phenethyl, 2,3,4-tris(trifluoromethyl)benzyl, 3,4,5-tris(trifluoromethyl)phenethyl, 2,5,6-tris(trifluoromethyl)benzyl, 2,4,6-tris(trifluoromethyl)phenethyl, 2,3,6-tris(tribromomethyl)benzyl, 2,3,4-tris(dibromomethyl)phenethyl, 3,4,5-tris(tribromomethyl)benzyl, 2,5,6-tris(dichloromethyl)phenethyl, 2,4,6-tris(dichloromethyl)benzyl, 1-trifluoromethyl-2-naphthylethyl, 2-trifluoromethyl-1-naphthylmethyl, 3-trifluoromethyl-1-naphthylethyl, 1-trichloromethyl-2-naphthylmethyl, 2-dichloromethyl-1-naphthylethyl, 3-tribromomethyl-1-naphthylmethyl, 3,8-bis(trifluoromethyl)-1-naphthylethyl, 2,3-bis(trifluoromethyl)-1-naphthylmethyl, 4,8-bis(trifluoromethyl)-1-naphthylethyl, 5,6-bis(trifluoromethyl)-1-naphthylmethyl, 3,8-bis(trichloromethyl)-1-naphthylethyl, 2,3-bis(dichloromethyl)-1-naphthylmethyl, 4,8-bis(dibromomethyl)-1-naphthylethyl, 5,6-bis(tribromomethyl)-1-naphthylmethyl, 2,3,6-tris(trifluoromethyl)-1-naphthylethyl, 2,3,4-tris(trifluoromethyl)-1-naphthylmethyl, 3,4,5-tris(trifluoromethyl)-1-naphthylethyl, 4,5,6-tris(trifluoromethyl)-1-naphthylmethyl, 2,4,8-tris(trifluoromethyl)-1-naphthylmethyl, bis(4-trifluoromethylphenyl)methyl, 4-trifluoromethylphenylphenylmethyl, bis(2-trichloromethylphenyl)methyl, bis(3-trichloromethylphenyl)methyl, bis(4-trichloromethylphenyl)methyl, 2-tribromomethylphenylphenylmethyl, 3-tribromomethylphenylphenylmethyl, bis(4-tribromomethylphenyl)methyl, bis(3,5-bis(trifluoromethyl)phenyl)methyl, bis(2,5-bis(trifluoromethyl)phenyl)methyl, bis(2,6-bis(trifluoromethyl)phenyl)methyl, 2,4-bis(trifluoromethyl)phenylphenylmethyl, bis(3,5-bis(tribromomethyl)phenyl)methyl, 2,5-bis(tribromomethyl)phenylphenylmethyl, 2,6-bis(trichloromethyl)phenylphenylmethyl, bis(2,4-bis(trichloromethyl)phenyl)methyl and bis(2,3,6-tris(trifluoromethyl)phenyl)methyl groups; aralkyl groups substituted with at least one lower alkyl group (as defined above), examples of which include 2-methylbenzyl, 3-methylbenzyl, 4-methylbenzyl, 2-methylphenethyl, 4-methylphenethyl, 2-ethylbenzyl, 3-propylphenethyl, 4-ethylbenzyl, 2-butylphenethyl, 3-pentylbenzyl, 4-pentylphenethyl, 3,5-dimethylbenzyl, 2,5-dimethylphenethyl, 2,6-dimethylbenzyl, 2,4-dimethylphenethyl, 3,5-dibutylbenzyl, 2,5-dipentylphenethyl, 2,6-dipropylbenzyl, 2,4-dipropylphenethyl, 2,3,6-trimethylbenzyl, 2,3,4-trimethylphenethyl, 3,4,5-trimethylbenzyl, 2,4,6-trimethylbenzyl, 2,5,6-trimethylphenethyl, 2,3,6-tributylphenethyl, 2,3,4-tripentylbenzyl, 3,4,5-tributylphenethyl, 2,5,6-tripropylbenzyl, 2,4,6-tripropylphenethyl, 1-methyl-2-naphthylmethyl, 2-methyl-1-naphthylethyl, 3-methyl-1-naphthylmethyl, 1-ethyl-2-naphthylethyl, 2-propyl-1-naphthylmethyl, 3-butyl-1-naphthylethyl, 3,8-dimethyl-1-naphthylmethyl, 2,3-dimethyl-1-naphthylethyl, 4,3-dimethyl-1-naphthylmethyl, 5,6-dimethyl-1-naphthylethyl, 3,8-diethyl-1-naphthylmethyl, 2,3-dipropyl-1-naphthylmethyl, 4,8-dipentyl-1-naphthylethyl, 5,6-dibutyl-1-naphthylmethyl, 2,3,6-trimethyl-1-naphthylmethyl, 2,3,4-trimethyl-1-naphthylethyl, 3,4,5-trimethyl-1-naphthylmethyl, 4,5,6-trimethyl-1-naphthylmethyl, 2,4,8-trimethyl-1-naphthylmethyl, bis(2-methylphenyl)methyl, 3-methylphenylphenylmethyl, bis(4-methylphenyl)methyl, 4-methylphenylphenylmethyl, bis(2-ethylphenyl)methyl, bis(3-ethylphenyl)methyl, bis(4-ethylphenyl)methyl, 2-propylphenylphenylmethyl, 3-propylphenylphenylmethyl, bis(4-propylphenyl)methyl, bis(3,5-dimethylphenyl)methyl, bis(2,5-dimethylphenyl)methyl, bis(2,6-dimethylphenyl)methyl, 2,4-dimethylphenylphenylmethyl, bis(3,5-dipropylphenyl)methyl, 2,5-dipropylphenylphenylmethyl, 2,6-diethylphenylphenylmethyl, bis(2,4-diethylphenyl)methyl and bis(2,3,6-trimethylphenyl)methyl groups; aralkyl groups substituted with at least one lower alkoxy group (as defined below), examples of which include 2-methoxybenzyl, 3-methoxybenzyl, 4-methoxybenzyl, 3-methoxyphenethyl, 2-ethoxyphenethyl, 3-propoxybenzyl, 4-ethoxyphenethyl, 2-butoxybenzyl, 3-pentoxyphenethyl, 4-pentoxybenzyl, 3,5-dimethoxyphenethyl, 2,5-dimethoxybenzyl, 2,6-dimethoxyphenethyl, 2,4-dimethoxybenzyl, 3,5-dibutoxyphenethyl, 2,5-dipentoxybenzyl, 2,6-dipropoxyphenethyl, 2,4-dipropoxybenzyl, 2,3,6-trimethoxyphenethyl, 2,3,4-trimethoxybenzyl, 3,4,5-trimethoxyphenethyl, 2,5,6-trimethoxybenzyl, 2,4,6-trimethoxyphenethyl, 2,3,6-tributoxybenzyl, 2,3,4-tripentoxyphenethyl, 3,4,5-tributoxybenzyl, 2,5,6-tripropoxyphenethyl, 2,4,6-tripropoxybenzyl, 1-methoxy-2-naphthylmethyl, 2-methoxy-1-naphthylmethyl, 3-methoxy-1-naphthylethyl, 1-ethoxy-2-naphthylmethyl, 2-propoxy-1-naphthylmethyl, 3-butoxy-1-naphthylethyl, 3,3-dimethoxy-1-naphthylmethyl, 2,3-dimethoxy-1-naphthylmethyl, 4,8-dimethoxy-1-naphthylethyl, 5,6-dimethoxy-1-naphthylmethyl, 3,8-diethoxy-1-naphthylmethyl, 2,3-dipropoxy-1-naphthylethyl, 4,8-dipentoxy-1-naphthylmethyl, 5,6-dibutoxy-1-naphthylethyl, 2,3,6-trimethoxy-1-naphthylethyl, 2,3,4-trimethoxy-1-naphthylmethyl, 3,4,5-trimethoxy-1-naphthylmethyl, 4,5,6-trimethoxy-1-naphthylethyl, 2,4,8-trimethoxy-1-naphthylmethyl, bis(2-methoxyphenyl)methyl, 3-methoxyphenylphenylmethyl, bis(4-methoxyphenyl)methyl, 4-methoxyphenylphenylmethyl, bis(2-ethoxyphenyl)methyl, bis(3-ethoxyphenyl)methyl, bis(4-ethoxyphenyl)methyl, 2-propoxyphenylphenylmethyl, 3-propoxyphenylphenylmethyl, bis(4-propoxyphenyl)methyl, bis(3,5-dimethoxyphenyl)methyl, bis(2,5-dimethoxyphenyl)methyl, bis(2,6-dimethoxyphenyl)methyl, 2,4-dimethoxyphenylphenylmethyl, bis(3,5-dipropoxyphenyl)methyl, 2,5-dipropoxyphenylphenylmethyl, 2,6-diethoxyphenylphenylmethyl, bis(2,4-diethoxyphenyl)methyl and bis(2,3,6-trimethoxyphenyl)methyl groups; aralkyl groups substituted with at least one amino group, examples of which include 2-aminophenethyl, 3-aminobenzyl, 4-aminophenethyl, 3,5-diaminobenzyl, 2,5-diaminophenethyl, 2,6-diaminobenzyl, 2,4-diaminophenethyl, 2,3,6-triaminobenzyl, 2,3,4-triaminophenethyl, 3,4,5-triaminobenzyl, 2,5,6-triaminophenethyl, 2,4,6-triaminobenzyl, 1-amino-2-naphthylmethyl, 2-amino-1-naphthylethyl, 3-amino-1-naphthylmethyl, 3,8-diamino-1-naphthylmethyl, 2,3-diamino-1-naphthylethyl, 4,8-diamino-1-naphthylmethyl, 5,6-diamino-1-naphthylmethyl, 2,3,6-triamino-1-naphthylethyl, 2,3,4-triamino-1-naphthylmethyl, 3,4,5-triamino-1-naphthylmethyl, 4,5,6-triamino-1-naphthylethyl, 2,4,8-triamino-1-naphthylmethyl, bis(2-aminophenyl)methyl, 3-amino-phenylphenylmethyl, bis(4-aminophenyl)methyl, 4-aminophenylphenylmethyl, bis(3,5-diaminophenyl)methyl, bis(2,5-diaminophenyl)methyl, bis(2,6-diaminophenyl)methyl, 2,4-diaminophenylphenylmethyl and bis(2,3,6-triaminophenyl)methyl groups; aralkyl groups substituted with at least

one nitro group, examples of which include 2-nitrophenethyl, 3-nitrobenzyl, 4-nitrobenzyl, 4-nitrophenethyl, 3,5-dinitrobenzyl, 2,5-dinitrophenethyl, 2,6-dinitrobenzyl, 2,4-dinitrophenethyl, 2,3,6-trinitrobenzyl, 2,3,4-trinitrophenethyl, 3,4,5-trinitrobenzyl, 2,5,6-trinitrophenethyl, 2,4,6-trinitrobenzyl, 1-nitro-2-naphthylmethyl, 2-nitro-1-naphthylethyl, 3-nitro-1-naphthylmethyl, 3,8-dinitro-1-naphthylmethyl, 2,3-dinitro-1-naphthylethyl, 4,8-dinitro-1-naphthylmethyl, 5,6-dinitro-1-naphthylmethyl, 2,3,6-trinitro-1-naphthylethyl, 2,3,4-trinitro-1-naphthylmethyl, 3,4,5-trinitro-1-naphthylmethyl, 4,5,6-trinitro-1-naphthylethyl, 2,4,8-trinitro-1-naphthylmethyl, bis(2-nitrophenyl)methyl, 3-nitrophenylphenylmethyl, bis(4-nitrophenyl)methyl, 4-nitrophenylphenylmethyl, bis(3,5-dinitrophenyl)methyl, bis(2,5-dinitrophenyl)methyl, bis(2,6-dinitrophenyl)methyl, 2,4-dinitrophenylphenylmethyl and bis(2,3,6-trinitrophenyl)methyl groups; and aralkyl groups substituted with at least one cyano group, examples of which include 2-cyanophenethyl, 3-cyanobenzyl, 4-cyanobenzyl, 4-cyanobenzylidiphenylmethyl, 4-cyanophenethyl, 3,5-dicyanobenzyl, 2,5-dicyanophenethyl, 2,6-dicyanobenzyl, 2,4-dicyanophenethyl, 2,3,6-tricyanobenzyl, 2,3,4-tricyanophenethyl, 3,4,5-tricyanobenzyl, 2,5,6-tricyanophenethyl, 2,4,6-tricyanobenzyl, 1-cyano-2-naphthylmethyl, 3-cyano-1-naphthylmethyl, 3,8-dicyano-1-naphthylmethyl, 2,3-dicyano-1-naphthylethyl, 4,8-dicyano-1-naphthylmethyl, 5,6-dicyano-1-naphthylmethyl, 2,3,6-tricyano-1-naphthylethyl, 2,3,4-tricyano-1-naphthylmethyl, 3,4,5-tricyano-1-naphthylmethyl, 4,5,6-tricyano-1-naphthylethyl, 2,4,8-tricyano-1-naphthylmethyl, bis(2-cyanophenyl)methyl, 3-cyanophenylphenylmethyl, bis(4-cyanophenyl)methyl, 4-cyanophenylphenylmethyl, bis(3,5-dicyanophenyl)methyl, bis(2,5-dicyanophenyl)methyl, bis(2,6-dicyanophenyl)methyl, 2,4-dicyanophenylphenylmethyl and bis(2,3,6-tricyanophenyl)methyl groups.

[0051] Of the above, unsubstituted aralkyl groups and aralkyl groups substituted with at least one substituent selected from the group consisting of halogen atoms, lower alkyl groups and lower alkoxy groups are preferred, unsubstituted aralkyl groups and aralkyl groups substituted with at least one substituent selected from the group consisting of halogen atoms and lower alkyl groups are more preferred, and unsubstituted aralkyl groups are most preferred.

[0052] Where substituents R^a , R^b , R^c or R^d represent a lower alkylsulfonyl group, this is a group in which a lower alkyl group, defined and exemplified above, is bonded to a sulfonyl group ($-SO_2-$). The lower alkylsulfonyl group is preferably a straight or branched alkylsulfonyl group having from 1 to 4 carbon atoms, more preferably a methylsulfonyl, ethylsulfonyl, propylsulfonyl, isopropylsulfonyl or butylsulfonyl group, and most preferably a methylsulfonyl, ethylsulfonyl or propylsulfonyl group.

[0053] Where substituents R^c and R^d together with the nitrogen atom to which they are bonded form a heterocyclyl group, said heterocyclyl group is a heterocyclyl group having at least one nitrogen atom as defined above in relation to substituent R^d . Examples of said heterocyclyl group include 1-azetidiny, 1-pyrrolidiny, 1-pyrroliny, 1-imidazolidiny, 1-imidazoliny, 1-pyrazolidiny, 1-pyrazoliny, 3-oxazolidiny, 3-thiazolidiny, 1-piperidy, tetrahydropyridin-1-yl, dihydropyridin-1-yl, 1-piperaziny, 4-morpholiny, 4-thiomorpholiny, 1-homopiperidy, 8-azabicyclo[3.2.1]octan-8-yl, 8-azabicyclo[3.2.1]octen-8-yl, 9-azabicyclo[3.3.1]nonan-9-yl and 9-azabicyclo[3.3.1]nonen-9-yl groups.

[0054] Where substituents R^c and R^d together with the nitrogen atom to which they are bonded form a heterocyclyl group as defined and exemplified above, said heterocyclyl groups may be fused with another cyclic group selected from the group consisting of aryl groups defined above and heteroaryl groups defined above. Examples of such fused heterocyclyl groups include tetrahydroquinolin-1-yl and tetrahydroisoquinolin-2-yl.

[0055] The halogen atoms in the definition of Substituent group α include fluorine, chlorine, bromine and iodine atoms, of which fluorine and chlorine atoms are preferred.

[0056] Where the substituent in the definition of Substituent group α is a lower alkoxy group, this is a group in which an oxygen atom is bonded to a lower alkyl group as defined and exemplified above. The alkoxy groups are preferably straight or branched alkoxy groups having 1 to 4 carbon atoms, more preferably methoxy, ethoxy, propoxy, isopropoxy or butoxy groups, and particularly preferably methoxy, ethoxy or propoxy groups.

[0057] Where the substituent in the definition of Substituent group α is a halogeno lower alkoxy group this is a group in which a lower alkoxy group as defined above is substituted with at least one halogen atom as exemplified above. The halogeno lower alkoxy groups preferably have from 1 to 4 carbon atoms, and are more preferably selected from the group consisting of difluoromethoxy, trifluoromethoxy and 2,2,2-trifluoromethoxy groups. Difluoromethoxy groups are most preferred.

[0058] Where the substituent in the definition of Substituent group α is a lower alkylthio group this is a group in which a sulfur atom is bonded to a lower alkyl group as defined and exemplified above. The lower alkylthio groups are preferably straight or branched alkylthio groups having 1 to 4 carbon atoms, more preferably methylthio, ethylthio, propylthio, isopropylthio or butylthio groups, and particularly preferably methylthio, ethylthio or propylthio groups.

[0059] Where the substituent in the definition of Substituent group α is a halogeno lower alkylthio group this is a group in which a lower alkylthio group as defined above is substituted with at least one halogen atom as exemplified above. The halogeno lower alkylthio groups preferably have from 1 to 4 carbon atoms, and are more preferably selected from the group consisting of difluoromethylthio, trifluoromethylthio and 2,2,2-trifluoroethylthio groups.

[0060] A preferred group of substituents of Substituent group α is Substituent group α^1 which comprises halogen atoms, lower alkoxy groups as defined above and halogeno lower alkoxy groups as defined above.

[0061] A preferred group of substituents of Substituent group β is Substituent group β^1 which comprises a group of

formula $-NR^cR^d$ wherein one of R^c and R^d is selected from the group consisting of hydrogen atoms and lower alkyl groups as defined above and the other is selected from the group consisting of hydrogen atoms, lower alkyl groups as defined above and aralkyl groups as defined above.

[0062] A preferred group of substituents of Substituent group γ is Substituent group γ^1 which comprises lower alkyl groups as defined above which are substituted with a substituent selected from the group consisting of amino groups, amino groups substituted with one or two lower alkyl groups as defined above and amino groups substituted with an aralkyl group as defined above. Of these, substituents in which the alkyl moiety which is substituted with a substituent selected from the group consisting of amino groups, amino groups substituted with one or two lower alkyl groups and amino groups substituted with an aralkyl is an alkyl group having from 1 to 4 carbon atoms are preferred. Aminomethyl, 2-aminoethyl, 3-aminopropyl, methylaminomethyl, 2-(methylamino)ethyl, 3-(methylamino)propyl, ethylaminomethyl, 2-(ethylamino)ethyl, 3-(ethylamino)propyl, dimethylaminomethyl, 2-(dimethylamino)ethyl, 3-(dimethylamino)propyl, diethylaminomethyl, 2-(diethylamino)ethyl, 3-(diethylamino)propyl, benzylaminomethyl, 2-(benzylamino)ethyl and 3-(benzylamino)propyl groups are more preferred.

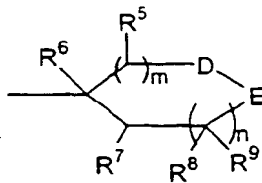
[0063] A preferred group of substituents of Substituent group δ is Substituent group δ^1 which comprises lower alkyl groups as defined above, halogeno lower alkyl groups, hydroxy lower alkyl groups and nitro lower alkyl groups.

[0064] The halogeno lower alkyl groups in the definition of Substituent group δ^1 above are lower alkyl groups as defined above which are substituted with at least one halogen atom. Halogeno alkyl groups having from 1 to 4 carbon atoms are preferred, trifluoromethyl, trichloromethyl, difluoromethyl, dichloromethyl, dibromomethyl, fluoromethyl, 2,2,2-trichloroethyl, 2,2,2-trifluoroethyl, 2-bromoethyl, 2-chloroethyl, 2-fluoroethyl and 2,2-dibromoethyl groups are more preferred, trifluoromethyl, trichloromethyl, difluoromethyl and fluoromethyl groups are still more preferred, and trifluoromethyl groups are most preferred.

[0065] The hydroxy lower alkyl groups in the definition of Substituent group δ^1 above are lower alkyl groups as defined above which are substituted with at least one hydroxy group. Hydroxyalkyl groups having from 1 to 4 carbon atoms are preferred, and hydroxymethyl, 2-hydroxyethyl and 3-hydroxypropyl groups are most preferred.

[0066] The nitro lower alkyl groups in the definition of Substituent group δ^1 above are lower alkyl groups as defined above which are substituted with at least one nitro group. Nitroalkyl groups having from 1 to 4 carbon atoms are preferred, and nitromethyl, 2-nitroethyl and 3-nitropropyl are most preferred.

[0067] Where R^3 is a group of formula $-X-R^4$ wherein X is a single bond and R^4 is a heterocyclyl group which may be optionally substituted with at least one group selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which is substituted with at least one group selected from the group consisting of Substituent group β defined above and Substituent group γ defined above or it is a heterocyclyl group having at least one nitrogen atom which may be optionally substituted with at least one group selected from the group consisting of Substituent group α defined above and Substituent group δ defined above, the group R^3 is preferably a group of the following formula (II):



wherein:

m represents 0 or 1;

n represents 1 or 2;

R^5 , R^6 , R^7 , R^8 and R^9 may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

one of D and E represents a group of formula $>NR^{10}$ wherein R^{10} is selected from the group consisting of hydrogen atoms, Substituent group γ defined above and

Substituent group δ defined above, and the other represents a group of formula $>CR^{11}R^{12}$ wherein R^{11} and R^{12} are the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above, or

R⁶ may form, together with R⁵ or R⁷, a single bond, and/or R¹⁰ and R¹¹ together may form a lower alkylene group defined below which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above.

[0068] Preferably, one of R⁵ and R⁷ forms a single bond together with R⁶, and the other is selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , more preferably selected from the group consisting of hydrogen atoms, lower alkyl groups as defined above and aralkyl groups as defined above, and most preferably selected from the group consisting of hydrogen atoms, alkyl groups having from 1 to 4 carbon atoms, benzyl groups and phenethyl groups.

[0069] Each of R⁸, R⁹, R¹⁰, R¹¹ and R¹² is preferably selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , more preferably selected from the group consisting of hydrogen atoms, lower alkyl groups as defined above and aralkyl groups as defined above, and most preferably selected from the group consisting of hydrogen atoms, alkyl groups having from 1 to 4 carbon atoms, benzyl groups and phenethyl groups.

[0070] Where R¹⁰ and R¹¹ together form a lower alkylene group, this is a straight or branched alkylene group having from 1 to 6 carbon atoms, examples of which include methylene, ethylene, trimethylene, propylene, tetramethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethylethylene, pentamethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 1,2-dimethyltrimethylene and hexamethylene groups. Straight or branched alkylene groups having from 3 to 6 carbon atoms are preferred, straight or branched alkylene group having 3 or 4 carbon atoms are more preferred, and straight chain alkylene groups having 3 or 4 carbon atoms are most preferred.

[0071] Preferred examples of the above group of formula (II) include:

(a) a group of formula (II) wherein m is 1 and n is 1;

(b) a group of formula (II) wherein one of R⁵ and R⁷ together with R⁶ represents a single bond and the other is a hydrogen atom;

(c) a group of formula (II) wherein R⁷, R⁸ and R⁹ may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group γ defined above and Substituent group δ defined above;

(d) a group of formula (II) wherein R¹⁰, R¹¹ and R¹² may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group γ defined above and Substituent group δ defined above;

(e) a group of formula (II) wherein R¹⁰ and R¹¹ together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above; and

(f) a group of formula (II) wherein each of R⁷, R⁸ and R⁹ is a hydrogen atom, and R¹⁰ and R¹¹ together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above.

[0072] Of these, more preferred examples of the above group of formula (II) include:

(g) a group of formula (II) wherein m is 1, n is 1, D represents a group of formula >CR¹¹R¹² and E represents a group of formula >NR¹⁰; and

(h) a group of formula (II) wherein m is 1, n is 1, D represents a group of formula >CR¹¹R¹², E represents a group of formula >NR¹⁰, and at least one of the substituents R⁸, R⁹, R¹¹ and R¹² is a substituent selected from the group consisting of Substituent group α defined above, Substituent group γ defined above and Substituent group δ defined above.

[0073] We also prefer a group of formula -X-R⁴ wherein X is a lower alkenylene group which may optionally be substituted with at least one substituent selected from Substituent group α defined above and R⁴ is a heterocyclyl group which may optionally be substituted with at least one group selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which is substituted with at least one group selected from the group consisting of Substituent group β defined above and Substituent group γ defined above or it is a heterocyclyl group having at least one nitrogen atom which may optionally be substituted with at least one group selected from the group consisting of Substituent group α defined above and Substituent group δ defined above.

[0074] In this case, preferred examples of such a heterocyclyl group for R⁴ include 2-pyrrolidinyl, 1-methyl-2-pyrro-

lidinyl, 1-ethyl-2-pyrrolidinyl, 1-propyl-2-pyrrolidinyl, 2-piperidinyl, 1-methyl-2-piperidinyl, 1-ethyl-2-piperidinyl and 1-propyl-2-piperidinyl groups.

[0075] The present invention encompasses esters and other derivatives of the compounds of formula (I). These esters and other derivatives are compounds of formula (I) in which a functional group (for example, a hydroxyl group, an amino group, an imino group or a sulfonamide group) of said compound of formula (I) is modified by the addition of a protecting group using conventional techniques well-known in the art (see, for example, "Protective Groups in Organic Synthesis, Second Edition, Theodora W. Greene and Peter G.M. Wuts, 1991, John Wiley & Sons, Inc.).

[0076] There is no particular restriction on the nature of this protecting group, provided that, where the ester or other derivative is intended for therapeutic purposes, it must be pharmacologically acceptable, i.e. the protecting group must be capable of being removed by a metabolic process (e.g. hydrolysis) on administration of said compound to the body of a live mammal to give a compound of formula (I) or a salt thereof. In other words, the pharmacologically acceptable esters or other derivatives are pro-drugs of the compounds of formula (I) of the present invention. Where, however, the ester or other derivative of the compound of formula (I) of the present invention is intended for non-therapeutic purposes (e.g. as an intermediate in the preparation of other compounds), then the requirement that said ester or other derivative is pharmacologically acceptable does not apply.

[0077] Whether an ester or other derivative of a compound of formula (I) of the present invention is pharmacologically acceptable can be easily determined. The compound under investigation is intravenously administered to an experimental animal such as a rat or mouse and the body fluids of the animal are thereafter studied. If a compound of formula (I) or a pharmacologically acceptable salt thereof can be detected in the body fluids, the compound under investigation is judged to be a pharmacologically acceptable ester or other derivative.

[0078] The compounds of formula (I) of the present invention can be converted to an ester, examples of which include a compound of formula (I) in which a hydroxyl group present therein is esterified. The ester residue may be a general protecting group where the esterified compound is to be used as an intermediate or a protecting group which is capable of being removed by a metabolic process (e.g. hydrolysis) *in vivo* where the esterified compound is one which is pharmacologically acceptable.

[0079] The general protecting group referred to above is a protecting group which is removable by a chemical process such as hydrolysis, hydrogenolysis, electrolysis or photolysis. Preferred examples of such a general protecting group used to synthesise a compound of formula (I) in which a hydroxyl residue therein is esterified include the following:

(i) aliphatic acyl groups, examples of which include

alkylcarbonyl groups having from 1 to 25 carbon atoms, examples of which include formyl, acetyl, propionyl, butyryl, isobutyryl, pentanoyl, pivaloyl, valeryl, isovaleryl, octanoyl, nonanoyl, decanoyl, 3-methylnonanoyl, 8-methylnonanoyl, 3-ethyloctanoyl, 3,7-dimethyloctanoyl, undecanoyl, dodecanoyl, tridecanoyl, tetradecanoyl, pentadecanoyl, hexadecanoyl, 1-methylpentadecanoyl, 14-methylpentadecanoyl, 13,13-dimethyltetradecanoyl, heptadecanoyl, 15-methylhexadecanoyl, octadecanoyl, 1-methylheptadecanoyl, nonadecanoyl, eicosanoyl and heneicosanoyl groups,

halogenated alkylcarbonyl groups having from 1 to 25 carbons in which the alkyl moiety thereof is substituted by at least one halogen atom, examples of which include chloroacetyl, dichloroacetyl, trichloroacetyl and trifluoroacetyl groups,

lower alkoxyalkylcarbonyl groups which comprise an alkylcarbonyl group having from 1 to 25 carbon atoms in which the alkyl moiety thereof is substituted with at least one lower alkoxy group as defined above, examples of said lower alkoxyalkylcarbonyl groups including methoxyacetyl groups, and

unsaturated alkylcarbonyl groups having from 1 to 25 carbon atoms, examples of which include acryloyl, propioloyl, methacryloyl, crotonoyl, isocrotonoyl and (E)-2-methyl-2-butenoyl groups; of these, alkylcarbonyl groups having from 1 to 6 carbon atoms are preferred;

(ii) aromatic acyl groups, examples of which include

arylcabonyl groups which comprise a carbonyl group which is substituted with an aryl group as defined above, examples of which include benzoyl, α -naphthoyl and β -naphthoyl groups,

halogenated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with at least one halogen atom, examples of which include 2-bromobenzoyl and 4-chlorobenzoyl groups,

lower alkylated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with at least one lower alkyl group as defined above, examples of which include 2,4,6-trimethyl-benzoyl and 4-toluoyl groups,

lower alkoxyated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with at least one lower alkoxy group as defined above, examples of which include 4-anisoyl groups,

nitroated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with at least one nitro group, examples of which include 4-nitrobenzoyl and 2-nitrobenzoyl groups;

lower alkoxy-carbonylated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with a carbonyl group which is itself substituted with a lower alkoxy group as defined above,

examples of which include 2-(methoxycarbonyl)benzoyl groups, and

arylated arylcarbonyl groups which comprise an arylcarbonyl group as defined above which is substituted with at least one aryl group as defined above, examples of which include 4-phenylbenzoyl groups;

(iii) alkoxy-carbonyl groups, examples of which include

lower alkoxy-carbonyl groups which comprise a carbonyl group substituted with a lower alkoxy group as defined above, examples of which include methoxycarbonyl, ethoxycarbonyl, propoxycarbonyl, butoxycarbonyl, s-butoxycarbonyl, t-butoxycarbonyl and isobutoxycarbonyl groups, and

lower alkoxy-carbonyl groups as defined above which are substituted with at least one substituent selected from the group consisting of halogen atoms and tri(lower alkyl)silyl groups (wherein said lower alkyl groups are as defined above), examples of which include 2,2,2-trichloroethoxycarbonyl and 2-trimethylsilylethoxycarbonyl groups;

(iv) tetrahydropyranyl or tetrahydrothiopyranyl groups which may optionally be substituted with at least one substituent selected from lower alkyl groups as defined above, halogen atoms and lower alkoxy groups as defined above, examples of which include tetrahydropyran-2-yl, 3-bromotetrahydropyran-2-yl, 4-methoxytetrahydropyran-4-yl, tetrahydrothiopyran-2-yl and 4-methoxytetrahydrothiopyran-4-yl groups;

(v) tetrahydrofuranyl or tetrahydrothiofuranyl groups which may optionally be substituted with at least one substituent selected from lower alkyl groups as defined above, halogen atoms and lower alkoxy groups as defined above, examples of which include tetrahydrofuran-2-yl and tetrahydrothiofuran-2-yl groups;

(vi) silyl groups, examples of which include

tri(lower alkyl)silyl groups (wherein said lower alkyl groups are as defined above), examples of which include trimethylsilyl, triethylsilyl, isopropyl dimethylsilyl, t-butyl dimethylsilyl, methyl diisopropylsilyl, methyl-di-t-butylsilyl and triisopropylsilyl groups, and

tri(lower alkyl)silyl groups in which at least one of said lower alkyl groups is substituted with 1 or 2 aryl groups as defined above, examples of which include diphenylmethylsilyl, diphenylbutylsilyl, diphenylisopropylsilyl and phenyl diisopropylsilyl groups;

(vii) alkoxy-methyl groups, examples of which include

lower alkoxy-methyl groups which comprise a methyl group which is substituted with a lower alkoxy group as defined above, examples of which include methoxymethyl, 1,1-dimethyl-1-methoxymethyl, ethoxymethyl, propoxymethyl, isopropoxymethyl, butoxymethyl and t-butoxymethyl groups,

lower alkoxy-lated lower alkoxy-methyl groups which comprise a lower alkoxy-methyl group as defined above in which the alkoxy moiety thereof is substituted with a lower alkoxy group as defined above, examples of which include 2-methoxyethoxymethyl groups, and

lower halogeno alkoxy-methyl groups which comprise a lower alkoxy-methyl group as defined above in which the alkoxy moiety thereof is substituted with at least one halogen atom, examples of which include 2,2,2-trichloroethoxymethyl and bis(2-chloroethoxy)methyl groups;

(viii) substituted ethyl groups, examples of which include

lower alkoxy-lated ethyl groups which comprise an ethyl group which is substituted with a lower alkoxy group as defined above, examples of which include 1-ethoxyethyl and 1-(isopropoxy)ethyl groups, and halogenated ethyl groups such as 2,2,2-trichloroethyl groups;

(ix) aralkyl groups as defined above, examples of which include

lower alkyl groups as defined above which are substituted with from 1 to 3 aryl groups as defined above, examples of which include benzyl, α -naphthylmethyl, β -naphthylmethyl, diphenylmethyl, triphenylmethyl, α -

naphthyldiphenylmethyl and 9-anthrylmethyl groups, and

lower alkyl groups as defined above which are substituted with from 1 to 3 aryl groups as defined above in which said aryl moiety is substituted with at least one substituent selected from the group consisting of lower alkyl groups as defined above, lower alkoxy groups as defined above, nitro groups, halogen atoms and cyano groups, examples of which include 4-methylbenzyl, 2,4,6-trimethylbenzyl, 3,4,5-trimethylbenzyl, 4-methoxybenzyl, 4-methoxyphenyldiphenylmethyl, 2-nitrobenzyl, 4-nitrobenzyl, 4-chlorobenzyl, 4-bromobenzyl and 4-cyanobenzyl groups;

(x) "alkenyloxycarbonyl groups" which comprise a carbonyl group which is substituted with an alkenyloxy group having from 2 to 6 carbon atoms, examples of which include vinyloxycarbonyl and allyloxycarbonyl groups; and

(xi) aralkyloxycarbonyl groups which comprise a carbonyl group which is substituted with an aralkyloxy group (which is an oxygen atom substituted with an aralkyl group as defined above), in which the aryl moiety thereof may optionally be substituted with one or two substituents selected from lower alkoxy groups as defined above and nitro groups, examples of which include benzyloxycarbonyl, 4-methoxybenzyloxycarbonyl, 3,4-dimethoxybenzyloxycarbonyl, 2-nitrobenzyloxycarbonyl and 4-nitrobenzyloxycarbonyl groups.

[0080] The protecting group which is capable of being removed by a metabolic process (e.g. hydrolysis) *in vivo* is one, which on administration to the body of a live mammal is removable by a metabolic process (e.g. hydrolysis) to give a compound of formula (I) or a salt thereof. Preferred examples of such a protecting group which is used to synthesise a compound of formula (I) in which a hydroxyl residue therein is esterified include the following:

(i) 1-(acyloxy)lower alkyl groups, examples of which include

1-(aliphatic acyloxy)lower alkyl groups which comprise a lower alkyl group as defined above which is substituted with an alkylcarbonyloxy group having from 1 to 6 carbon atoms, examples of which include formyloxymethyl, acetoxymethyl, dimethylaminoacetoxymethyl, propionyloxymethyl, butyryloxymethyl, pivaloyloxymethyl, valeryloxymethyl, isovaleryloxymethyl, hexanoyloxymethyl, 1-formyloxyethyl, 1-acetoxyethyl, 1-propionyloxyethyl, 1-butyryloxyethyl, 1-pivaloyloxyethyl, 1-valeryloxyethyl, 1-isovaleryloxyethyl, 1-hexanoyloxyethyl, 1-formyloxypropyl, 1-acetoxypropyl, 1-propionyloxypropyl, 1-butyryloxypropyl, 1-pivaloyloxypropyl, 1-valeryloxypropyl, 1-isovaleryloxypropyl, 1-hexanoyloxypropyl, 1-acetoxybutyl, 1-propionyloxybutyl, 1-butyryloxybutyl, 1-pivaloyloxybutyl, 1-acetoxypentyl, 1-propionyloxypentyl, 1-butyryloxypentyl, 1-pivaloyloxypentyl and 1-pivaloyloxyhexyl groups,

1-(cycloalkylcarbonyloxy)lower alkyl groups which comprise a lower alkyl group as defined above which is substituted with a cycloalkylcarbonyloxy group in which a carbonyloxy group is substituted with a cyclohexyl group as defined above, examples of which include cyclopentylcarbonyloxymethyl, cyclohexylcarbonyloxymethyl, 1-cyclopentylcarbonyloxyethyl, 1-cyclohexylcarbonyloxyethyl, 1-cyclopentylcarbonyloxypropyl, 1-cyclohexylcarbonyloxypropyl, 1-cyclopentylcarbonyloxybutyl and 1-cyclohexylcarbonyloxybutyl groups, and

1-(aromatic acyloxy)lower alkyl groups which comprise a lower alkyl group as defined above which is substituted with an aralkyloxycarbonyl group which comprises a carbonyl group which is substituted with an aralkyloxy group (said aralkyloxy group comprising an oxygen atom which is substituted with an aralkyl group as defined above), examples of which include benzyloxymethyl groups;

(ii) substituted carbonyloxyalkyl groups, examples of which include

(lower alkoxy carbonyloxy)alkyl groups which comprise a lower alkyl group as defined above which is substituted with a lower alkoxy carbonyloxy group which comprises a carbonyloxy group substituted with a lower alkoxy group as defined above, examples of which include methoxycarbonyloxymethyl, ethoxycarbonyloxymethyl, propoxycarbonyloxymethyl, isopropoxycarbonyloxymethyl, butoxycarbonyloxymethyl, isobutoxycarbonyloxymethyl, pentyloxycarbonyloxymethyl, hexyloxycarbonyloxymethyl, cyclohexyloxycarbonyloxymethyl, cyclohexyloxycarbonyloxy(cyclohexyl)methyl, 1-(methoxycarbonyloxy)ethyl, 1-(ethoxycarbonyloxy)ethyl, 1-(propoxycarbonyloxy)ethyl, 1-(isopropoxycarbonyloxy)ethyl, 1-(butoxycarbonyloxy)ethyl, 1-(isobutoxycarbonyloxy)ethyl, 1-(tert-butoxycarbonyloxy)ethyl, 1-(pentyloxycarbonyloxy)ethyl, 1-(hexyloxycarbonyloxy)ethyl, 1-(cyclopentyloxycarbonyloxy)ethyl, 1-(cyclohexyloxycarbonyloxy)propyl, 1-(cyclopentyloxycarbonyloxy)propyl, 1-(cyclohexyloxycarbonyloxy)propyl, 1-(ethoxycarbonyloxy)propyl, 1-(propoxycarbonyloxy)propyl, 1-(isopropoxycarbonyloxy)propyl, 1-(butoxycarbonyloxy)propyl, 1-(isobutoxycarbonyloxy)propyl, 1-(pentyloxycarbonyloxy)propyl, 1-(hexyloxycarbonyloxy)propyl, 1-(methoxycarbonyloxy)butyl, 1-(ethoxycarbonyloxy)butyl, 1-(propoxycarbonyloxy)butyl, 1-(isopropoxycarbonyloxy)butyl, 1-(butoxycarbonyloxy)butyl, 1-(isobutoxycarbonyloxy)butyl, 1-(methoxy-

carbonyloxy)pentyl, 1-(ethoxycarbonyloxy)pentyl, 1-(methoxycarbonyloxy)hexyl and 1-(ethoxycarbonyloxy)hexyl groups, and oxodioxolenylmethyl groups, which comprise a methyl group which is substituted with an oxodioxolenyl group which itself may optionally be substituted with a group selected from the group consisting of lower alkyl groups as defined above and aryl groups as defined above which may optionally be substituted with at least one lower alkyl group as defined above, lower alkoxy group as defined above or halogen atom, examples of which include (5-phenyl-2-oxo-1,3-dioxolen-4-yl)methyl, [5-(4-methylphenyl)-2-oxo-1,3-dioxolen-4-yl]methyl, [5-(4-methoxyphenyl)-2-oxo-1,3-dioxolen-4-yl]methyl, [5-(4-fluorophenyl)-2-oxo-1,3-dioxolen-4-yl]methyl, [5-(4-chlorophenyl)-2-oxo-1,3-dioxolen-4-yl]methyl, (2-oxo-1,3-dioxolen-4-yl)-methyl, (5-methyl-2-oxo-1,3-dioxolen-4-yl)methyl, (5-ethyl-2-oxo-1,3-dioxolen-4-yl)methyl, (5-propyl-2-oxo-1,3-dioxolen-4-yl)methyl, (5-isopropyl-2-oxo-1,3-dioxolen-4-yl)methyl and (5-butyl-2-oxo-1,3-dioxolen-4-yl)methyl groups;

(iii) phthalidyl groups which comprise a phthalidyl group which may optionally be substituted with a substituent selected from the group consisting of lower alkyl groups as defined above and lower alkoxy groups as defined above, examples of which include phthalidyl, dimethylphthalidyl and dimethoxyphthalidyl groups;

(iv) aliphatic acyl groups as defined and exemplified above in relation to the general protecting group for a hydroxyl group

(v) aromatic acyl groups as defined and exemplified above in relation to the general protecting group for a hydroxyl group

(vi) half-ester salt residues of succinic acid;

(vii) phosphate ester salt residues;

(viii) ester-forming residues of an amino acid;

(ix) carbamoyl groups which may optionally be substituted with 1 or 2 lower alkyl groups as defined above; and

(x) 1-(acyloxy)alkoxycarbonyl groups which comprise a lower alkoxycarbonyl group as defined above in which the lower alkoxy moiety is substituted with an aliphatic acyloxy group as defined above or an aromatic acyloxy group as defined above, examples of which include pivaloyloxymethyloxycarbonyl groups.

[0081] Of the above protecting groups which are capable of being removed by a metabolic process (e.g. hydrolysis) *in vivo* which are used to synthesise a compound of formula (I) in which a hydroxyl residue therein is esterified, the substituted carbonyloxyalkyl groups are preferred.

[0082] In the case where the compound of formula (I) of the present invention has an amino group, an imino group and/or a sulfonamide group, the compound can be converted to a derivative other than the esters described above and the pharmacologically acceptable salts described below. The "other derivatives" of the compounds of formula (I) include such derivatives. Example of such derivatives include an amide derivative in which an aliphatic acyl group defined and exemplified above or an aromatic acyl group defined and exemplified above is bonded to a nitrogen atom of an amino group, imino group and/or sulfonamide group present in said compound of formula (I). Where said derivative is a pharmacologically acceptable derivative of a compound of formula (I) it must be capable of being removed by a metabolic process (e.g. hydrolysis) on administration of said compound to the body of a live mammal to give a compound of formula (I) or a salt thereof.

[0083] Where the compound of formula (I) of the present invention or a pharmacologically acceptable ester or other derivative thereof has a basic group, such as an amino group, the compound can be converted to a salt by reacting it with an acid, and in the case where the compound of formula (I) of the present invention or a pharmacologically acceptable ester or other derivative thereof has an acidic group, such as a sulfonamide group, the compound can be converted to a salt by reacting it with a base. The compounds of the present invention encompass such salts. Where said salts are to be used for a therapeutic use, they must be pharmacologically acceptable.

[0084] Preferred examples of the salts formed with a basic group present in the compound of formula (I) of the present invention include inorganic acid salts such as hydrohalogenated acid salts (e.g. hydrochlorides, hydrobromides and hydroiodides), nitrates, perchlorates, sulfates and phosphates; organic acid salts such as lower alkanesulfonates in which the lower alkyl moiety thereof is as defined above (e.g. methanesulfonates, trifluoromethanesulfonates and ethanesulfonates), arylsulfonates in which the aryl moiety thereof is as defined above (e.g. benzenesulfonate or p-toluenesulfonate), acetates, malates, fumarates, succinates, citrates, ascorbates, tartrates, oxalates and maleates; and amino acid salts such as glycine salts, lysine salts, arginine salts, ornithine salts, glutamates and aspartates.

[0085] Preferred example of the salts formed with an acidic group present in the compound of formula (I) of the present invention include metal salts such as alkali metal salts (e.g. sodium salts, potassium salts and lithium salts), alkali earth metal salts (e.g. calcium salts and magnesium salts), aluminum salts and iron salts; amine salts such as inorganic amine salts (e.g. ammonium salts) and organic amine salts (e.g. t-octylamine salts, dibenzylamine salts, morpholine salts, glucosamine salts, phenylglycinealkyl ester salts, ethylenediamine salts, N-methylglucamine salts, guanidine salts, diethylamine salts, triethylamine salts, dicyclohexylamine salts, N,N'-dibenzylethylenediamine salts, chloroprocaine salts, procaine salts, diethanolamine salts, N-benzylphenethylamine salts, piperazine salts, tetramethylammonium salts and tris(hydroxymethyl)aminomethane salts; and amino acid salts such as glycine salts, lysine salts, arginine salts, ornithine salts, glutamates and aspartates.

[0086] The compound of formula (I) of the present invention can sometimes take up water upon exposure to the atmosphere or when recrystallized to absorb water or to form a hydrate and such hydrates are also included within the scope of the present invention. Additionally, certain other solvents may be taken up by the compounds of the present invention to produce solvates, which also form a part of the present invention.

[0087] The compounds of formula (I) of the present invention can sometimes exist in the form of geometrical isomers (cis and trans isomers) and, where said compounds contain one or more asymmetric centres, optical isomers. For the compounds of the present invention, each of said isomers and mixture of said isomers are depicted by a single formula, i.e. the formula (I). Accordingly, the present invention covers both the individual isomers and mixtures thereof in any proportion, including racemic mixtures.

[0088] Preferred classes of compounds of the present invention are those compounds of formula (I) and pharmacologically acceptable salts, esters and other derivatives thereof wherein:

(A) R¹ is an aryl group defined above which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(B) R¹ is a phenyl or naphthyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(C) R¹ is a phenyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α^1 defined above, Substituent group β^1 defined above, Substituent group γ^1 defined above and Substituent group δ^1 defined above;

(D) R¹ is a phenyl group which may optionally be substituted with at least one substituent selected from the group consisting of halogen atoms, halogeno lower alkyl groups as defined above and a halogeno lower alkoxy groups as defined above;

(E) R¹ is a phenyl, 4-fluorophenyl, 3-fluorophenyl, 3-chlorophenyl, 3,4-difluorophenyl, 3,4,5-trifluorophenyl, 3-chloro-4-fluorophenyl, 3-difluoromethoxyphenyl or 3-trifluoromethylphenyl group;

(F) R² is a 5- or 6-membered aromatic heterocyclic group containing one or two nitrogen atoms, said group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(G) R² is a pyridyl or pyrimidinyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(H) R² is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(I) R² is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(J) R² is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a

substituent selected from the group consisting of methoxy, amino, methylamino, benzylamino, and α -methylbenzylamino groups;

(K) R^4 is selected from the group consisting of:

cycloalkyl groups having from 3 to 7 carbon atoms, which may be optionally substituted with at least one substituent selected from the group consisting of

Substituent group α defined above and Substituent group δ defined above and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ ;

phenyl or naphthyl groups which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above;

4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally one further heteroatom selected from the group consisting of oxygen atoms, sulfur atom and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above;

5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above;

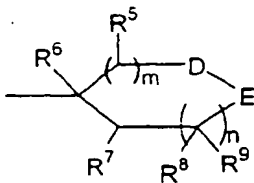
5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ ; and groups of formula $-NR^aR^b$ wherein R^a and R^b are as defined above;

(L) X is selected from the group consisting of single bonds, alkylene groups having from 1 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above, alkenylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above and alkynylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above;

(M) X is selected from the group consisting of single bonds, alkylene groups having from 1 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above, alkenylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above and alkynylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α defined above;

(N) X is a single bond or an alkenylene group having from 2 to 4 carbon atoms;

(O) R^3 is a group of the following formula (II):



wherein

m represents 0 or 1,

n represents 1 or 2,

R^5 , R^6 , R^7 , R^8 and R^9 may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

one of D and E represents a group of formula $>NR^{10}$ wherein R^{10} is selected from the group consisting of hydrogen atoms, Substituent group γ defined above and Substituent group δ defined above, and the other represents a group of formula $>CR^{11}R^{12}$ wherein R^{11} and R^{12} are the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above; or R^5 may form, together with R^6 or R^7 , a single bond, and/or R^{10} and R^{11} together may form a straight or branched alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(P) R^2 is a group of formula (II) as defined above wherein m is 1 and n is 1;

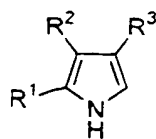
(Q) R^2 is a group of formula (II) as defined above wherein E represents a group of formula $>NR^{10}$ and D represents a group of formula $>CR^{11}R^{12}$;

(R) R^2 is a group of formula (II) as defined above wherein one of R^5 and R^7 together with R^6 represents a single bond and the other represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group γ defined above and Substituent group δ defined above, and R^8 , R^9 , R^{10} , R^{11} and R^{12} may be the same or different from one another and each represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α defined above, Substituent group γ defined above and Substituent group δ defined above, or R^{10} and R^{11} together form an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

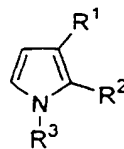
(S) R^2 is a group of formula (II) as defined above wherein R^{10} and R^{11} together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above;

(T) R^2 is a group of formula (II) as defined above wherein each of R^7 , R^8 and R^9 is a hydrogen atom, and R^{10} and R^{11} together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above; and

(U) said compound of formula (I) is a compound of formula (I-1) or (I-3) below wherein R^1 , R^2 and R^3 are as defined above;



(I-1)



(I-3)

[0089] Compounds of formula (I) which comprise any combination of the factors selected freely from the five groups consisting of (A) to (E) above; (F) to (J) above; (K) above; (L) to (N) above; (O) to (T) above; and (U) above are preferred.

[0090] More preferred compounds of the present invention are compounds of formula (I) and pharmacologically acceptable salts, esters and other derivatives thereof, wherein:

(i) R¹ is as defined in (A) above, R² is as defined in (F) above and R³ is a group of formula -X-R⁴ wherein X is as defined in (L) above and R⁴ is as defined in (K) above;

(ii) R¹ is as defined in (A) above, R² is as defined in (F) above and R³ is a group of formula -X-R⁴ wherein X is as defined in (M) above and R⁴ is as defined in (K) above;

(iii) R¹ is as defined in (B) above, R² is as defined in (G) above and R³ is a group of formula -X-R⁴ wherein X is as defined in (N) above and R⁴ is as defined in (K) above;

(iv) R¹ is as defined in (B) above, R² is as defined in (G) above and R³ is as defined in (O) above;

(v) R¹ is as defined in (C) above, R² is as defined in (H) above and R³ is as defined in (P) above;

(vi) R¹ is as defined in (C) above, R² is as defined in (H) above and R³ is as defined in (Q) above;

(vii) R¹ is as defined in (D) above, R² is as defined in (I) above and R³ is as defined in (R) above;

(viii) R¹ is as defined in (E) above, R² is as defined in (J) above and R³ is as defined in (S) above;

(ix) R¹ is as defined in (E) above, R² is as defined in (J) above and R³ is as defined in (T) above; and

(x) any of the above compounds (i) to (ix) wherein the compound of formula (I) is a compound of formula (I-1) or (I-3) as defined in (U) above.

[0091] Particularly preferred compounds of the present invention are compounds of formula (I) selected from the following group of compounds, and pharmacologically acceptable salts, esters and other derivatives thereof:

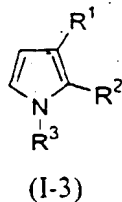
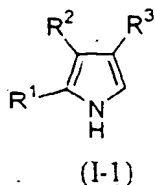
4-(3-aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(3-acetylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(3-methylaminopropyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 [5-(4-fluorophenyl)-4-(pyridin-4-yl)-1 *H*-pyrrol-3-yl]-(pyridin-4-yl)methanol,
 4-(1-acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(piperidin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 1-(1-acetylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-[1-(2-nitroethyl)piperidin-4-yl]-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-3-yl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 1-(azetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(3-dimethylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-[3-(piperidin-1-yl)propyl]-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-[3-(1-methylpiperazin-4-yl)propyl]-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(2,2,6,6-tetramethylpiperidin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-3-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(quinuclidin-3-yl)-1 *H*-pyrrole,
 1-(4-aminocyclohexyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1-(piperidin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydro-2,2,6,6-tetramethylpyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(2,2,6,6-tetramethylpiperidin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,

2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethylpiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,6-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,2,6,6-pentamethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(3-dimethylamino-1-propen-1-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(4-aminobutyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 1-(3-dimethylaminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,7,8,8a-octahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(6-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(2-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(6-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(2-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,9,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1 *H*-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1 *H*-pyrrole,
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1 *H*-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1 *H*-pyrrole, and
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1 *H*-pyrrole.

[0092] The most preferred compounds of the present invention are compounds of formula (I) selected from the following group of compounds, and pharmacologically acceptable salts, esters and other derivatives thereof:

2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,6-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,

2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1*H*-pyrrole. Specific examples of compounds of formula (I) of the present invention include the following compounds of formula (I-1) (Table 1) and formula (I-3) (Table 2).



In these Tables, the following abbreviations are used:

ABN	9-azabicyclo[3.3.1]nonanyl
dcH-ABN	dchydro[9-azabicyclo[3.3.1]nonanyl] (i.e. 9-azabicyclo[3.3.1]nonenyl)
ABO	8-azabicyclo[3.2.1]octanyl
dcH-ABO	dehydro[8-azabicyclo[3.2.1]octanyl] (i.e. 8-azabicyclo[3.2.1]octenyl)
Ac	acetyl
Azt	azetidiny
Bn	benzyl
Bu	butyl
t-Bu	t-butyl
Bz	benzoyl
Et	ethyl
Hp	heptyl
Hpip	homopiperidiny
Hx	hexyl
cHx	cyclohexyl
Ind	indoliziny
Me	methyl
Mor	morpholiny
Nn	nonyl
Oc	octyl
Ph	phenyl
Phet	phenethyl
Pip	piperidyl
deH-Pip	dehydropiperidyl (i.e. tetrahydropyridyl)
Piz	piperaziny
Pn	pentyl
cPn	cyclopentyl
Pr	propyl
iPr	isopropyl
cPr	cyclopropyl
Pym	pyrimidiny
Pyr	pyridyl
Pyrd	pyrrolidiny
Qui	quinoliziny
Qun	quinuclidiny

deH-Qun dehydroquinuclidinyl (i.e. quinuclidienyl)
Tmor thiomorpholinyl.

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Table 1

Compound No.	R ¹	R ²	R ³
1-1	Ph	4-PyT	H ₂ N-CH ₂
1-2	Ph	4-PyT	H ₂ N-(CH ₂) ₂
1-3	Ph	4-PyT	H ₂ N-(CH ₂) ₃
1-4	Ph	4-PyT	H ₂ N-(CH ₂) ₄
1-5	Ph	4-PyT	MeNH-CH ₂
1-6	Ph	4-PyT	MeNH-(CH ₂) ₂
1-7	Ph	4-PyT	MeNH-(CH ₂) ₃
1-8	Ph	4-PyT	MeNH-(CH ₂) ₄
1-9	Ph	4-PyT	EtNH-CH ₂
1-10	Ph	4-PyT	EtNH-(CH ₂) ₂
1-11	Ph	4-PyT	EtNH-(CH ₂) ₃
1-12	Ph	4-PyT	EtNH-(CH ₂) ₄
1-13	Ph	4-PyT	Me ₂ N-CH ₂
1-14	Ph	4-PyT	Me ₂ N-(CH ₂) ₂
1-15	Ph	4-PyT	Me ₂ N-(CH ₂) ₃
1-16	Ph	4-PyT	Me ₂ N-(CH ₂) ₄
1-17	Ph	4-PyT	1-Azt-CH ₂
1-18	Ph	4-PyT	1-Azt-(CH ₂) ₂
1-19	Ph	4-PyT	1-Azt-(CH ₂) ₃
1-20	Ph	4-PyT	1-Azt-(CH ₂) ₄
1-21	Ph	4-PyT	1-Pyrd-CH ₂
1-22	Ph	4-PyT	1-Pyrd-(CH ₂) ₂
1-23	Ph	4-PyT	1-Pyrd-(CH ₂) ₃
1-24	Ph	4-PyT	1-Pyrd-(CH ₂) ₄
1-25	Ph	4-PyT	1-Pip-CH ₂
1-26	Ph	4-PyT	1-Pip-(CH ₂) ₂
1-27	Ph	4-PyT	1-Pip-(CH ₂) ₃
1-28	Ph	4-PyT	1-Pip-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
1-29	Ph	4-Pyr	1-Mor-CH ₂
1-30	Ph	4-Pyr	1-Mor-(CH ₂) ₂
1-31	Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-32	Ph	4-Pyr	1-Mor-(CH ₂) ₄
1-33	Ph	4-Pyr	1-Tmor-CH ₂
1-34	Ph	4-Pyr	1-Tmor-(CH ₂) ₂
1-35	Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-36	Ph	4-Pyr	1-Tmor-(CH ₂) ₄
1-37	Ph	4-Pyr	1-Piz-CH ₂
1-38	Ph	4-Pyr	1-Piz-(CH ₂) ₂
1-39	Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-40	Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-41	Ph	4-Pyr	1-Piz-(CH ₂) ₄
1-42	Ph	4-Pyr	3-Azt
1-43	Ph	4-Pyr	1-Me-3-Azt
1-44	Ph	4-Pyr	1-Bn-3-Azt
1-45	Ph	4-Pyr	3-Pyrd
1-46	Ph	4-Pyr	1-Me-3-Pyrd
1-47	Ph	4-Pyr	3-Pip
1-48	Ph	4-Pyr	4-Pip
1-49	Ph	4-Pyr	4-(3,4-deH-Pip)
1-50	Ph	4-Pyr	1-Me-4-Pip
1-51	Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-52	Ph	4-Pyr	1-Et-4-Pip
1-53	Ph	4-Pyr	1-Bn-4-Pip
1-54	Ph	4-Pyr	3-Hpip
1-55	Ph	4-Pyr	1-Me-3-Hpip
1-56	Ph	4-Pyr	4-Hpip
1-57	Ph	4-Pyr	1-Me-4-Hpip
1-58	Ph	4-Pyr	2-Mor

Compound No.	R ¹	R ²	R ³
1-59	Ph	4-Pyr	1-Me-2-Mor
1-60	Ph	4-Pyr	2-Tmor
1-61	Ph	4-Pyr	1-Me-2-Tmor
1-62	Ph	4-Pyr	1-Piz
1-63	Ph	4-Pyr	4-Me-1-Piz
1-64	Ph	4-Pyr	2-Piz
1-65	Ph	4-Pyr	4-Pyr
1-66	Ph	4-Pyr	3-Pyr
1-67	Ph	4-Pyr	2-Pyr
1-68	Ph	4-Pyr	4-Pym
1-69	Ph	4-Pyr	5-Pym
1-70	Ph	4-Pyr	2-Pym
1-71	Ph	4-Pyr	3-Azt-CH ₂
1-72	Ph	4-Pyr	1-Me-3-Azt-CH ₂
1-73	Ph	4-Pyr	3-Pyrd-CH ₂
1-74	Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-75	Ph	4-Pyr	4-Pip-CH ₂
1-76	Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-77	Ph	4-Pyr	3-Hpip-CH ₂
1-78	Ph	4-Pyr	1-Me-3-Hpip-CH ₂
1-79	Ph	4-Pyr	4-Hpip-CH ₂
1-80	Ph	4-Pyr	1-Me-4-Hpip-CH ₂
1-81	Ph	4-Pyr	2-Mor-CH ₂
1-82	Ph	4-Pyr	1-Me-2-Mor-CH ₂
1-83	Ph	4-Pyr	2-Tmor-CH ₂
1-84	Ph	4-Pyr	1-Me-2-Tmor-CH ₂
1-85	Ph	4-Pyr	1-Piz-CH ₂
1-86	Ph	4-Pyr	4-Me-1-Piz-CH ₂
1-87	Ph	4-Pyr	2-Piz-CH ₂
1-88	Ph	4-Pyr	4-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-89	Ph	4-Pyr	3-Pyr-CH ₂
1-90	Ph	4-Pyr	2-Pyr-CH ₂
1-91	Ph	4-Pyr	4-Pym-CH ₂
1-92	Ph	4-Pyr	5-Pym-CH ₂
1-93	Ph	4-Pyr	2-Pym-CH ₂
1-94	Ph	4-Pyr	H ₂ N-CH ₂ CH=CH
1-95	4-F-Ph	4-Pyr	H ₂ N-CH ₂
1-96	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
1-97	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-98	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
1-99	4-F-Ph	4-Pyr	MeNH-CH ₂
1-100	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₂
1-101	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-102	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₄
1-103	4-F-Ph	4-Pyr	EtNH-CH ₂
1-104	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₂
1-105	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-106	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₄
1-107	4-F-Ph	4-Pyr	Me ₂ N-CH ₂
1-108	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
1-109	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-110	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
1-111	4-F-Ph	4-Pyr	1-Azt-CH ₂
1-112	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₂
1-113	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-114	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₄
1-115	4-F-Ph	4-Pyr	1-Pyrd-CH ₂
1-116	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
1-117	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-118	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
1-119	4-F-Ph	4-Pyr	1-Pip-CH ₂
1-120	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₂
1-121	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-122	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₄
1-123	4-F-Ph	4-Pyr	1-Mor-CH ₂
1-124	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₂
1-125	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-126	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₄
1-127	4-F-Ph	4-Pyr	1-Tmor-CH ₂
1-128	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
1-129	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-130	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
1-131	4-F-Ph	4-Pyr	1-Piz-CH ₂
1-132	4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₂
1-133	4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-134	4-F-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-135	4-F-Ph	4-Pyr	4-Bn-1-Piz-(CH ₂) ₃
1-136	4-F-Ph	4-Pyr	3-Azt
1-137	4-F-Ph	4-Pyr	1-Me-3-Azt
1-138	4-F-Ph	4-Pyr	1-Bn-3-Azt
1-139	4-F-Ph	4-Pyr	3-Pyrd
1-140	4-F-Ph	4-Pyr	1-Me-3-Pyrd
1-141	4-F-Ph	4-Pyr	3-Pip
1-142	4-F-Ph	4-Pyr	4-Pip
1-143	4-F-Ph	4-Pyr	4-(3,4-deH-Pip)
1-144	4-F-Ph	4-Pyr	1-Me-4-Pip
1-145	4-F-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-146	4-F-Ph	4-Pyr	1-Et-4-Pip
1-147	4-F-Ph	4-Pyr	1-Bn-4-Pip
1-148	4-F-Ph	4-Pyr	3-Hpip

Compound No.	R ¹	R ²	R ³
1-149	4-F-Ph	4-Pyr	1-Me-3-Hpip
1-150	4-F-Ph	4-Pyr	4-Hpip
1-151	4-F-Ph	4-Pyr	1-Me-4-Hpip
1-152	4-F-Ph	4-Pyr	2-Mor
1-153	4-F-Ph	4-Pyr	1-Me-2-Mor
1-154	4-F-Ph	4-Pyr	2-Tmor
1-155	4-F-Ph	4-Pyr	1-Me-2-Tmor
1-156	4-F-Ph	4-Pyr	1-Piz
1-157	4-F-Ph	4-Pyr	4-Me-1-Piz
1-158	4-F-Ph	4-Pyr	2-Piz
1-159	4-F-Ph	4-Pyr	4-Pyr
1-160	4-F-Ph	4-Pyr	3-Pyr
1-161	4-F-Ph	4-Pyr	2-Pyr
1-162	4-F-Ph	4-Pyr	4-Pym
1-163	4-F-Ph	4-Pyr	5-Pym
1-164	4-F-Ph	4-Pyr	2-Pym
1-165	4-F-Ph	4-Pyr	3-Azt-CH ₂
1-166	4-F-Ph	4-Pyr	1-Me-3-Azt-CH ₂
1-167	4-F-Ph	4-Pyr	3-Pyrd-CH ₂
1-168	4-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-169	4-F-Ph	4-Pyr	4-Pip-CH ₂
1-170	4-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-171	4-F-Ph	4-Pyr	3-Hpip-CH ₂
1-172	4-F-Ph	4-Pyr	1-Me-3-Hpip-CH ₂
1-173	4-F-Ph	4-Pyr	4-Hpip-CH ₂
1-174	4-F-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
1-175	4-F-Ph	4-Pyr	2-Mor-CH ₂
1-176	4-F-Ph	4-Pyr	1-Me-2-Mor-CH ₂
1-177	4-F-Ph	4-Pyr	2-Tmor-CH ₂
1-178	4-F-Ph	4-Pyr	1-Me-2-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
1-179	4-F-Ph	4-Pyr	1-Piz-CH ₂
1-180	4-F-Ph	4-Pyr	4-Me-1-Piz-CH ₂
1-181	4-F-Ph	4-Pyr	2-Piz-CH ₂
1-182	4-F-Ph	4-Pyr	4-Pyr-CH ₂
1-183	4-F-Ph	4-Pyr	3-Pyr-CH ₂
1-184	4-F-Ph	4-Pyr	2-Pyr-CH ₂
1-185	4-F-Ph	4-Pyr	4-Pym-CH ₂
1-186	4-F-Ph	4-Pyr	5-Pym-CH ₂
1-187	4-F-Ph	4-Pyr	2-Pym-CH ₂
1-188	3-F-Ph	4-Pyr	H ₂ N-CH ₂
1-189	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
1-190	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-191	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
1-192	3-F-Ph	4-Pyr	MeNH-CH ₂
1-193	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₂
1-194	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-195	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₄
1-196	3-F-Ph	4-Pyr	EtNH-CH ₂
1-197	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₂
1-198	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-199	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₄
1-200	3-F-Ph	4-Pyr	Me ₂ N-CH ₂
1-201	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
1-202	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-203	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
1-204	3-F-Ph	4-Pyr	1-Azt-CH ₂
1-205	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₂
1-206	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-207	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₄
1-208	3-F-Ph	4-Pyr	1-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-209	3-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
1-210	3-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-211	3-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
1-212	3-F-Ph	4-Pyr	1-Pip-CH ₂
1-213	3-F-Ph	4-Pyr	1-Pip-(CH ₂) ₂
1-214	3-F-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-215	3-F-Ph	4-Pyr	1-Pip-(CH ₂) ₄
1-216	3-F-Ph	4-Pyr	1-Mor-CH ₂
1-217	3-F-Ph	4-Pyr	1-Mor-(CH ₂) ₂
1-218	3-F-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-219	3-F-Ph	4-Pyr	1-Mor-(CH ₂) ₄
1-220	3-F-Ph	4-Pyr	1-Tmor-CH ₂
1-221	3-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
1-222	3-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-223	3-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
1-224	3-F-Ph	4-Pyr	1-Piz-CH ₂
1-225	3-F-Ph	4-Pyr	1-Piz-(CH ₂) ₂
1-226	3-F-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-227	3-F-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-228	3-F-Ph	4-Pyr	1-Piz-(CH ₂) ₄
1-229	3-F-Ph	4-Pyr	3-Azt
1-230	3-F-Ph	4-Pyr	1-Me-3-Azt
1-231	3-F-Ph	4-Pyr	1-Bn-3-Azt
1-232	3-F-Ph	4-Pyr	3-Pyrd
1-233	3-F-Ph	4-Pyr	1-Me-3-Pyrd
1-234	3-F-Ph	4-Pyr	3-Pip
1-235	3-F-Ph	4-Pyr	4-Pip
1-236	3-F-Ph	4-Pyr	4-(3,4-deH-Pip)
1-237	3-F-Ph	4-Pyr	1-Me-4-Pip
1-238	3-F-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-239	3-F-Ph	4-Pyr	1-Et-4-Pip
1-240	3-F-Ph	4-Pyr	1-Bn-4-Pip
1-241	3-F-Ph	4-Pyr	3-Hpip
1-242	3-F-Ph	4-Pyr	1-Me-3-Hpip
1-243	3-F-Ph	4-Pyr	4-Hpip
1-244	3-F-Ph	4-Pyr	1-Me-4-Hpip
1-245	3-F-Ph	4-Pyr	2-Mor
1-246	3-F-Ph	4-Pyr	1-Me-2-Mor
1-247	3-F-Ph	4-Pyr	2-Tmor
1-248	3-F-Ph	4-Pyr	1-Me-2-Tmor
1-249	3-F-Ph	4-Pyr	1-Piz
1-250	3-F-Ph	4-Pyr	4-Me-1-Piz
1-251	3-F-Ph	4-Pyr	2-Piz
1-252	3-F-Ph	4-Pyr	4-Pyr
1-253	3-F-Ph	4-Pyr	3-Pyr
1-254	3-F-Ph	4-Pyr	2-Pyr
1-255	3-F-Ph	4-Pyr	4-Pym
1-256	3-F-Ph	4-Pyr	5-Pym
1-257	3-F-Ph	4-Pyr	2-Pym
1-258	3-F-Ph	4-Pyr	3-Azt-CH ₂
1-259	3-F-Ph	4-Pyr	1-Me-3-Azt-CH ₂
1-260	3-F-Ph	4-Pyr	3-Pyrd-CH ₂
1-261	3-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-262	3-F-Ph	4-Pyr	4-Pip-CH ₂
1-263	3-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-264	3-F-Ph	4-Pyr	3-Hpip-CH ₂
1-265	3-F-Ph	4-Pyr	1-Me-3-Hpip-CH ₂
1-266	3-F-Ph	4-Pyr	4-Hpip-CH ₂
1-267	3-F-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
1-268	3-F-Ph	4-Pyr	2-Mor-CH ₂

Compound No.	R ¹	R ²	R ³
1-269	3-F-Ph	4-Pyr	1-Me-2-Mor-CH ₂
1-270	3-F-Ph	4-Pyr	2-Tmor-CH ₂
1-271	3-F-Ph	4-Pyr	1-Me-2-Tmor-CH ₂
1-272	3-F-Ph	4-Pyr	1-Piz-CH ₂
1-273	3-F-Ph	4-Pyr	4-Me-1-Piz-CH ₂
1-274	3-F-Ph	4-Pyr	2-Piz-CH ₂
1-275	3-F-Ph	4-Pyr	4-Pyr-CH ₂
1-276	3-F-Ph	4-Pyr	3-Pyr-CH ₂
1-277	3-F-Ph	4-Pyr	2-Pyr-CH ₂
1-278	3-F-Ph	4-Pyr	4-Pym-CH ₂
1-279	3-F-Ph	4-Pyr	5-Pym-CH ₂
1-280	3-F-Ph	4-Pyr	2-Pym-CH ₂
1-281	3,4-diF-Ph	4-Pyr	H ₂ N-CH ₂
1-282	3,4-diF-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
1-283	3,4-diF-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-284	3,4-diF-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
1-285	3,4-diF-Ph	4-Pyr	MeNH-CH ₂
1-286	3,4-diF-Ph	4-Pyr	MeNH-(CH ₂) ₂
1-287	3,4-diF-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-288	3,4-diF-Ph	4-Pyr	MeNH-(CH ₂) ₄
1-289	3,4-diF-Ph	4-Pyr	EtNH-CH ₂
1-290	3,4-diF-Ph	4-Pyr	EtNH-(CH ₂) ₂
1-291	3,4-diF-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-292	3,4-diF-Ph	4-Pyr	EtNH-(CH ₂) ₄
1-293	3,4-diF-Ph	4-Pyr	Me ₂ N-CH ₂
1-294	3,4-diF-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
1-295	3,4-diF-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-296	3,4-diF-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
1-297	3,4-diF-Ph	4-Pyr	1-Azt-CH ₂
1-298	3,4-diF-Ph	4-Pyr	1-Azt-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
1-299	3,4-diF-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-300	3,4-diF-Ph	4-Pyr	1-Azt-(CH ₂) ₄
1-301	3,4-diF-Ph	4-Pyr	1-Pyrd-CH ₂
1-302	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
1-303	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-304	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
1-305	3,4-diF-Ph	4-Pyr	1-Pip-CH ₂
1-306	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₂
1-307	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-308	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₄
1-309	3,4-diF-Ph	4-Pyr	1-Mor-CH ₂
1-310	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₂
1-311	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-312	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₄
1-313	3,4-diF-Ph	4-Pyr	1-Tmor-CH ₂
1-314	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
1-315	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-316	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
1-317	3,4-diF-Ph	4-Pyr	1-Piz-CH ₂
1-318	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₂
1-319	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-320	3,4-diF-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-321	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₄
1-322	3,4-diF-Ph	4-Pyr	3-Azt
1-323	3,4-diF-Ph	4-Pyr	1-Me-3-Azt
1-324	3,4-diF-Ph	4-Pyr	1-Bn-3-Azt
1-325	3,4-diF-Ph	4-Pyr	3-Pyrd
1-326	3,4-diF-Ph	4-Pyr	1-Me-3-Pyrd
1-327	3,4-diF-Ph	4-Pyr	3-Pip
1-328	3,4-diF-Ph	4-Pyr	4-Pip

Compound No.	R ¹	R ²	R ³
1-329	3,4-diF-Ph	4-Pyr	4-(3,4-deH-Pip)
1-330	3,4-diF-Ph	4-Pyr	1-Me-4-Pip
1-331	3,4-diF-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-332	3,4-diF-Ph	4-Pyr	1-Et-4-Pip
1-333	3,4-diF-Ph	4-Pyr	1-Bn-4-Pip
1-334	3,4-diF-Ph	4-Pyr	3-Hpip
1-335	3,4-diF-Ph	4-Pyr	1-Me-3-Hpip
1-336	3,4-diF-Ph	4-Pyr	4-Hpip
1-337	3,4-diF-Ph	4-Pyr	1-Me-4-Hpip
1-338	3,4-diF-Ph	4-Pyr	2-Mor
1-339	3,4-diF-Ph	4-Pyr	1-Me-2-Mor
1-340	3,4-diF-Ph	4-Pyr	2-Tmor
1-341	3,4-diF-Ph	4-Pyr	1-Me-2-Tmor
1-342	3,4-diF-Ph	4-Pyr	1-Piz
1-343	3,4-diF-Ph	4-Pyr	4-Me-1-Piz
1-344	3,4-diF-Ph	4-Pyr	2-Piz
1-345	3,4-diF-Ph	4-Pyr	4-Pyr
1-346	3,4-diF-Ph	4-Pyr	3-Pyr
1-347	3,4-diF-Ph	4-Pyr	2-Pyr
1-348	3,4-diF-Ph	4-Pyr	4-Pym
1-349	3,4-diF-Ph	4-Pyr	5-Pym
1-350	3,4-diF-Ph	4-Pyr	2-Pym
1-351	3,4-diF-Ph	4-Pyr	3-Azt-CH ₂
1-352	3,4-diF-Ph	4-Pyr	1-Me-3-Azt-CH ₂
1-353	3,4-diF-Ph	4-Pyr	3-Pyrd-CH ₂
1-354	3,4-diF-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-355	3,4-diF-Ph	4-Pyr	4-Pip-CH ₂
1-356	3,4-diF-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-357	3,4-diF-Ph	4-Pyr	3-Hpip-CH ₂
1-358	3,4-diF-Ph	4-Pyr	1-Me-3-Hpip-CH ₂

Compound No.	R ¹	R ²	R ³
1-359	3,4-diF-Ph	4-Pyr	4-Hpip-CH ₂
1-360	3,4-diF-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
1-361	3,4-diF-Ph	4-Pyr	2-Mor-CH ₂
1-362	3,4-diF-Ph	4-Pyr	1-Me-2-Mor-CH ₂
1-363	3,4-diF-Ph	4-Pyr	2-Tmor-CH ₂
1-364	3,4-diF-Ph	4-Pyr	1-Me-2-Tmor-CH ₂
1-365	3,4-diF-Ph	4-Pyr	1-Piz-CH ₂
1-366	3,4-diF-Ph	4-Pyr	4-Me-1-Piz-CH ₂
1-367	3,4-diF-Ph	4-Pyr	2-Piz-CH ₂
1-368	3,4-diF-Ph	4-Pyr	4-Pyr-CH ₂
1-369	3,4-diF-Ph	4-Pyr	3-Pyr-CH ₂
1-370	3,4-diF-Ph	4-Pyr	2-Pyr-CH ₂
1-371	3,4-diF-Ph	4-Pyr	4-Pym-CH ₂
1-372	3,4-diF-Ph	4-Pyr	5-Pym-CH ₂
1-373	3,4-diF-Ph	4-Pyr	2-Pym-CH ₂
1-374	3-Cl-Ph	4-Pyr	H ₂ N-CH ₂
1-375	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
1-376	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-377	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
1-378	3-Cl-Ph	4-Pyr	MeNH-CH ₂
1-379	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₂
1-380	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-381	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₄
1-382	3-Cl-Ph	4-Pyr	EtNH-CH ₂
1-383	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₂
1-384	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-385	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₄
1-386	3-Cl-Ph	4-Pyr	Me ₂ N-CH ₂
1-387	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
1-388	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-389	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
1-390	3-Cl-Ph	4-Pyr	1-Azt-CH ₂
1-391	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₂
1-392	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-393	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₄
1-394	3-Cl-Ph	4-Pyr	1-Pyrd-CH ₂
1-395	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
1-396	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-397	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
1-398	3-Cl-Ph	4-Pyr	1-Pip-CH ₂
1-399	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₂
1-400	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-401	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₄
1-402	3-Cl-Ph	4-Pyr	1-Mor-CH ₂
1-403	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₂
1-404	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-405	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₄
1-406	3-Cl-Ph	4-Pyr	1-Tmor-CH ₂
1-407	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
1-408	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-409	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
1-410	3-Cl-Ph	4-Pyr	1-Piz-CH ₂
1-411	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₂
1-412	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-413	3-Cl-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-414	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₄
1-415	3-Cl-Ph	4-Pyr	3-Azt
1-416	3-Cl-Ph	4-Pyr	1-Me-3-Azt
1-417	3-Cl-Ph	4-Pyr	1-Bn-3-Azt
1-418	3-Cl-Ph	4-Pyr	3-Pyrd

Compound No.	R ¹	R ²	R ³
1-419	3-Cl-Ph	4-Pyr	1-Me-3-Pyrd
1-420	3-Cl-Ph	4-Pyr	3-Pip
1-421	3-Cl-Ph	4-Pyr	4-Pip
1-422	3-Cl-Ph	4-Pyr	4-(3,4-deH-Pip)
1-423	3-Cl-Ph	4-Pyr	1-Me-4-Pip
1-424	3-Cl-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-425	3-Cl-Ph	4-Pyr	1-Et-4-Pip
1-426	3-Cl-Ph	4-Pyr	1-Bn-4-Pip
1-427	3-Cl-Ph	4-Pyr	3-Hpip
1-428	3-Cl-Ph	4-Pyr	1-Me-3-Hpip
1-429	3-Cl-Ph	4-Pyr	4-Hpip
1-430	3-Cl-Ph	4-Pyr	1-Me-4-Hpip
1-431	3-Cl-Ph	4-Pyr	2-Mor
1-432	3-Cl-Ph	4-Pyr	1-Me-2-Mor
1-433	3-Cl-Ph	4-Pyr	2-Tmor
1-434	3-Cl-Ph	4-Pyr	1-Me-2-Tmor
1-435	3-Cl-Ph	4-Pyr	1-Piz
1-436	3-Cl-Ph	4-Pyr	4-Mc-1-Piz
1-437	3-Cl-Ph	4-Pyr	2-Piz
1-438	3-Cl-Ph	4-Pyr	4-Pyr
1-439	3-Cl-Ph	4-Pyr	3-Pyr
1-440	3-Cl-Ph	4-Pyr	2-Pyr
1-441	3-Cl-Ph	4-Pyr	4-Pym
1-442	3-Cl-Ph	4-Pyr	5-Pym
1-443	3-Cl-Ph	4-Pyr	2-Pym
1-444	3-Cl-Ph	4-Pyr	3-Azt-CH ₂
1-445	3-Cl-Ph	4-Pyr	1-Me-3-Azt-CH ₂
1-446	3-Cl-Ph	4-Pyr	3-Pyrd-CH ₂
1-447	3-Cl-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-448	3-Cl-Ph	4-Pyr	4-Pip-CH ₂

Compound No.	R ¹	R ²	R ³
1-449	3-Cl-Ph	4-PyT	1-Me-4-Pip-CH ₂
1-450	3-Cl-Ph	4-PyT	3-Hpip-CH ₂
1-451	3-Cl-Ph	4-PyT	1-Me-3-Hpip-CH ₂
1-452	3-Cl-Ph	4-PyT	4-Hpip-CH ₂
1-453	3-Cl-Ph	4-PyT	1-Me-4-Hpip-CH ₂
1-454	3-Cl-Ph	4-PyT	2-Mor-CH ₂
1-455	3-Cl-Ph	4-PyT	1-Me-2-Mor-CH ₂
1-456	3-Cl-Ph	4-PyT	2-Tmor-CH ₂
1-457	3-Cl-Ph	4-PyT	1-Me-2-Tmor-CH ₂
1-458	3-Cl-Ph	4-PyT	1-Piz-CH ₂
1-459	3-Cl-Ph	4-PyT	4-Me-1-Piz-CH ₂
1-460	3-Cl-Ph	4-PyT	2-Piz-CH ₂
1-461	3-Cl-Ph	4-PyT	4-PyT-CH ₂
1-462	3-Cl-Ph	4-PyT	3-PyT-CH ₂
1-463	3-Cl-Ph	4-PyT	2-PyT-CH ₂
1-464	3-Cl-Ph	4-PyT	4-Pym-CH ₂
1-465	3-Cl-Ph	4-PyT	5-Pym-CH ₂
1-466	3-Cl-Ph	4-PyT	2-Pym-CH ₂
1-467	Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
1-468	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
1-469	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-470	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
1-471	Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
1-472	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
1-473	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-474	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
1-475	Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
1-476	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
1-477	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-478	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
1-479	Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
1-480	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
1-481	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-482	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
1-483	Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
1-484	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
1-485	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-486	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
1-487	Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
1-488	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
1-489	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-490	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
1-491	Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
1-492	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
1-493	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-494	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
1-495	Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
1-496	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
1-497	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-498	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
1-499	Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
1-500	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
1-501	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-502	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
1-503	Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-504	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-505	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-506	Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-507	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
1-508	Ph	2-NH ₂ -4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
1-509	Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-510	Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
1-511	Ph	2-NH ₂ -4-Pym	3-Pyrd
1-512	Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-513	Ph	2-NH ₂ -4-Pym	3-Pip
1-514	Ph	2-NH ₂ -4-Pym	4-Pip
1-515	Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-516	Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-517	Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-518	Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
1-519	Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
1-520	Ph	2-NH ₂ -4-Pym	3-Hpip
1-521	Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
1-522	Ph	2-NH ₂ -4-Pym	4-Hpip
1-523	Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
1-524	Ph	2-NH ₂ -4-Pym	2-Mor
1-525	Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
1-526	Ph	2-NH ₂ -4-Pym	2-Tmor
1-527	Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
1-528	Ph	2-NH ₂ -4-Pym	1-Piz
1-529	Ph	2-NH ₂ -4-Pym	4-Mc-1-Piz
1-530	Ph	2-NH ₂ -4-Pym	2-Piz
1-531	Ph	2-NH ₂ -4-Pym	4-Pyr
1-532	Ph	2-NH ₂ -4-Pym	3-Pyr
1-533	Ph	2-NH ₂ -4-Pym	2-Pyr
1-534	Ph	2-NH ₂ -4-Pym	4-Pym
1-535	Ph	2-NH ₂ -4-Pym	5-Pym
1-536	Ph	2-NH ₂ -4-Pym	2-Pym
1-537	Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
1-538	Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂

Compound No.	R ¹	R ²	R ³
1-539	Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-540	Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-541	Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-542	Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-543	Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
1-544	Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
1-545	Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
1-546	Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
1-547	Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
1-548	Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
1-549	Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
1-550	Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
1-551	Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-552	Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
1-553	Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-554	Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-555	Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-556	Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
1-557	Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-558	Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-559	Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-560	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
1-561	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
1-562	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-563	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
1-564	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
1-565	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
1-566	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-567	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
1-568	4-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂

Compound No.	R ¹	R ²	R ³
1-569	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
1-570	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-571	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
1-572	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
1-573	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
1-574	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-575	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
1-576	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
1-577	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
1-578	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-579	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
1-580	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
1-581	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
1-582	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-583	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
1-584	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
1-585	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
1-586	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-587	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
1-588	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
1-589	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
1-590	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-591	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
1-592	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
1-593	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
1-594	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-595	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
1-596	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-597	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-598	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-599	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-600	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-601	4-F-Ph	2-NH ₂ -4-Pym	3-Azt
1-602	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-603	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
1-604	4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-605	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-606	4-F-Ph	2-NH ₂ -4-Pym	3-Pip
1-607	4-F-Ph	2-NH ₂ -4-Pym	4-Pip
1-608	4-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-609	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-610	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-611	4-F-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
1-612	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
1-613	4-F-Ph	2-NH ₂ -4-Pym	3-Hpip
1-614	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
1-615	4-F-Ph	2-NH ₂ -4-Pym	4-Hpip
1-616	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
1-617	4-F-Ph	2-NH ₂ -4-Pym	2-Mor
1-618	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
1-619	4-F-Ph	2-NH ₂ -4-Pym	2-Tmor
1-620	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
1-621	4-F-Ph	2-NH ₂ -4-Pym	1-Piz
1-622	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-623	4-F-Ph	2-NH ₂ -4-Pym	2-Piz
1-624	4-F-Ph	2-NH ₂ -4-Pym	4-Pyr
1-625	4-F-Ph	2-NH ₂ -4-Pym	3-Pyr
1-626	4-F-Ph	2-NH ₂ -4-Pym	2-Pyr
1-627	4-F-Ph	2-NH ₂ -4-Pym	4-Pym
1-628	4-F-Ph	2-NH ₂ -4-Pym	5-Pym

Compound No.	R ¹	R ²	R ³
1-629	4-F-Ph	2-NH ₂ -4-Pym	2-Pym
1-630	4-F-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
1-631	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
1-632	4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-633	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-634	4-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-635	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-636	4-F-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
1-637	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
1-638	4-F-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
1-639	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
1-640	4-F-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
1-641	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
1-642	4-F-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
1-643	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
1-644	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-645	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
1-646	4-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-647	4-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-648	4-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-649	4-F-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
1-650	4-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-651	4-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-652	4-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-653	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
1-654	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
1-655	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-656	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
1-657	3-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
1-658	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
1-659	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-660	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
1-661	3-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
1-662	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
1-663	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-664	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
1-665	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
1-666	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
1-667	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-668	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
1-669	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
1-670	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
1-671	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-672	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
1-673	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
1-674	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
1-675	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-676	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
1-677	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
1-678	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
1-679	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-680	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
1-681	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
1-682	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
1-683	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-684	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
1-685	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
1-686	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
1-687	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-688	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
1-689	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-690	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-691	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-692	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₂
1-693	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
1-694	3-F-Ph	2-NH ₂ -4-Pym	3-Azt
1-695	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-696	3-F-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
1-697	3-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-698	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-699	3-F-Ph	2-NH ₂ -4-Pym	3-Pip
1-700	3-F-Ph	2-NH ₂ -4-Pym	4-Pip
1-701	3-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-702	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-703	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-704	3-F-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
1-705	3-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
1-706	3-F-Ph	2-NH ₂ -4-Pym	3-Hpip
1-707	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
1-708	3-F-Ph	2-NH ₂ -4-Pym	4-Hpip
1-709	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
1-710	3-F-Ph	2-NH ₂ -4-Pym	2-Mor
1-711	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
1-712	3-F-Ph	2-NH ₂ -4-Pym	2-Tmor
1-713	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
1-714	3-F-Ph	2-NH ₂ -4-Pym	1-Piz
1-715	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-716	3-F-Ph	2-NH ₂ -4-Pym	2-Piz
1-717	3-F-Ph	2-NH ₂ -4-Pym	4-Pyr
1-718	3-F-Ph	2-NH ₂ -4-Pym	3-Pyr

Compound No.	R ¹	R ²	R ³
1-719	3-F-Ph	2-NH ₂ -4-Pym	2-Pyr
1-720	3-F-Ph	2-NH ₂ -4-Pym	4-Pym
1-721	3-F-Ph	2-NH ₂ -4-Pym	5-Pym
1-722	3-F-Ph	2-NH ₂ -4-Pym	2-Pym
1-723	3-F-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
1-724	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
1-725	3-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-726	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-727	3-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-728	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-729	3-F-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
1-730	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
1-731	3-F-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
1-732	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
1-733	3-F-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
1-734	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
1-735	3-F-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
1-736	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
1-737	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-738	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
1-739	3-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-740	3-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-741	3-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-742	3-F-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
1-743	3-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-744	3-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-745	3-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-746	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
1-747	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
1-748	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-749	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
1-750	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
1-751	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
1-752	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-753	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
1-754	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
1-755	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
1-756	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-757	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
1-758	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
1-759	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
1-760	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-761	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
1-762	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
1-763	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
1-764	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-765	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
1-766	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
1-767	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
1-768	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-769	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
1-770	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
1-771	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
1-772	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-773	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
1-774	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
1-775	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
1-776	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-777	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
1-778	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
1-779	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
1-780	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-781	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
1-782	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-783	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-784	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-785	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-786	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
1-787	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Azt
1-788	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-789	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
1-790	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-791	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-792	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pip
1-793	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pip
1-794	3,4-diF-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-795	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-796	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-797	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
1-798	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
1-799	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Hpip
1-800	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
1-801	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Hpip
1-802	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
1-803	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Mor
1-804	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
1-805	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Tmor
1-806	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
1-807	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz
1-808	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz

Compound No.	R ¹	R ²	R ³
1-809	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Piz
1-810	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pyr
1-811	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyr
1-812	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pyr
1-813	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pym
1-814	3,4-diF-Ph	2-NH ₂ -4-Pym	5-Pym
1-815	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pym
1-816	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
1-817	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
1-818	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-819	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-820	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-821	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-822	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
1-823	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
1-824	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
1-825	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
1-826	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
1-827	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
1-828	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
1-829	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
1-830	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-831	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
1-832	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-833	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-834	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-835	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
1-836	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-837	3,4-diF-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-838	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂

Compound No.	R ¹	R ²	R ³
1-839	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
1-840	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
1-841	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-842	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
1-843	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
1-844	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
1-845	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-846	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
1-847	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
1-848	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
1-849	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-850	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
1-851	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
1-852	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
1-853	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-854	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
1-855	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
1-856	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
1-857	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-858	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
1-859	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
1-860	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
1-861	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-862	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
1-863	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
1-864	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
1-865	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-866	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
1-867	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
1-868	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
1-869	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-870	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
1-871	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
1-872	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
1-873	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-874	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
1-875	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-876	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
1-877	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-878	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-879	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
1-880	3-Cl-Ph	2-NH ₂ -4-Pym	3-Azt
1-881	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-882	3-Cl-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
1-883	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-884	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-885	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pip
1-886	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pip
1-887	3-Cl-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-888	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-889	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-890	3-Cl-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
1-891	3-Cl-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
1-892	3-Cl-Ph	2-NH ₂ -4-Pym	3-Hpip
1-893	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
1-894	3-Cl-Ph	2-NH ₂ -4-Pym	4-Hpip
1-895	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
1-896	3-Cl-Ph	2-NH ₂ -4-Pym	2-Mor
1-897	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
1-898	3-Cl-Ph	2-NH ₂ -4-Pym	2-Tmor

Compound No.	R ¹	R ²	R ³
1-899	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
1-900	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz
1-901	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-902	3-Cl-Ph	2-NH ₂ -4-Pym	2-Piz
1-903	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pyr
1-904	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyr
1-905	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pyr
1-906	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pym
1-907	3-Cl-Ph	2-NH ₂ -4-Pym	5-Pym
1-908	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pym
1-909	3-Cl-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
1-910	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
1-911	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-912	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-913	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-914	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-915	3-Cl-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
1-916	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
1-917	3-Cl-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
1-918	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
1-919	3-Cl-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
1-920	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
1-921	3-Cl-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
1-922	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
1-923	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
1-924	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
1-925	3-Cl-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-926	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-927	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-928	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-929	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-930	3-Cl-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-931	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-932	Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
1-933	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
1-934	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-935	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
1-936	Ph	2-MeNH-4-Pym	MeNH-CH ₂
1-937	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
1-938	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-939	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
1-940	Ph	2-MeNH-4-Pym	EtNH-CH ₂
1-941	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
1-942	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-943	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
1-944	Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
1-945	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
1-946	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-947	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
1-948	Ph	2-MeNH-4-Pym	1-Azt-CH ₂
1-949	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
1-950	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-951	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
1-952	Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
1-953	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
1-954	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-955	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
1-956	Ph	2-MeNH-4-Pym	1-Pip-CH ₂
1-957	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
1-958	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-959	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
1-960	Ph	2-MeNH-4-Pym	1-Mor-CH ₂
1-961	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
1-962	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-963	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
1-964	Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
1-965	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
1-966	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-967	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
1-968	Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-969	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
1-970	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-971	Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-972	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
1-973	Ph	2-MeNH-4-Pym	3-Azt
1-974	Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-975	Ph	2-MeNH-4-Pym	1-Bn-3-Azt
1-976	Ph	2-MeNH-4-Pym	3-Pyrd
1-977	Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-978	Ph	2-MeNH-4-Pym	3-Pip
1-979	Ph	2-MeNH-4-Pym	4-Pip
1-980	Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-981	Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-982	Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-983	Ph	2-MeNH-4-Pym	1-Et-4-Pip
1-984	Ph	2-MeNH-4-Pym	1-Bn-4-Pip
1-985	Ph	2-MeNH-4-Pym	3-Hpip
1-986	Ph	2-MeNH-4-Pym	1-Me-3-Hpip
1-987	Ph	2-MeNH-4-Pym	4-Hpip
1-988	Ph	2-MeNH-4-Pym	1-Me-4-Hpip

Compound No.	R ¹	R ²	R ³
1-989	Ph	2-MeNH-4-Pym	2-Mor
1-990	Ph	2-MeNH-4-Pym	1-Me-2-Mor
1-991	Ph	2-MeNH-4-Pym	2-Tmor
1-992	Ph	2-MeNH-4-Pym	1-Me-2-Tmor
1-993	Ph	2-MeNH-4-Pym	1-Piz
1-994	Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-995	Ph	2-MeNH-4-Pym	2-Piz
1-996	Ph	2-MeNH-4-Pym	4-Pyr
1-997	Ph	2-MeNH-4-Pym	3-Pyr
1-998	Ph	2-MeNH-4-Pym	2-Pyr
1-999	Ph	2-MeNH-4-Pym	4-Pym
1-1000	Ph	2-MeNH-4-Pym	5-Pym
1-1001	Ph	2-MeNH-4-Pym	2-Pym
1-1002	Ph	2-MeNH-4-Pym	3-Azt-CH ₂
1-1003	Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
1-1004	Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1005	Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1006	Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1007	Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1008	Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
1-1009	Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
1-1010	Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
1-1011	Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
1-1012	Ph	2-MeNH-4-Pym	2-Mor-CH ₂
1-1013	Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
1-1014	Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
1-1015	Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
1-1016	Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1017	Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
1-1018	Ph	2-MeNH-4-Pym	2-Piz-CH ₂

Compound No.	R ¹	R ²	R ³
1-1019	Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1020	Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1021	Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
1-1022	Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1023	Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1024	Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1025	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
1-1026	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
1-1027	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1028	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
1-1029	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂
1-1030	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
1-1031	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1032	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
1-1033	4-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂
1-1034	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
1-1035	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1036	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
1-1037	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
1-1038	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
1-1039	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1040	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
1-1041	4-F-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
1-1042	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
1-1043	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1044	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
1-1045	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
1-1046	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
1-1047	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1048	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
1-1049	4-F-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
1-1050	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
1-1051	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1052	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
1-1053	4-F-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
1-1054	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
1-1055	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1056	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
1-1057	4-F-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
1-1058	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
1-1059	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1060	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
1-1061	4-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1062	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
1-1063	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1064	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1065	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
1-1066	4-F-Ph	2-MeNH-4-Pym	3-Azt
1-1067	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1068	4-F-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
1-1069	4-F-Ph	2-MeNH-4-Pym	3-Pyrd
1-1070	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1071	4-F-Ph	2-MeNH-4-Pym	3-Pip
1-1072	4-F-Ph	2-MeNH-4-Pym	4-Pip
1-1073	4-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1074	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1075	4-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1076	4-F-Ph	2-MeNH-4-Pym	1-Et-4-Pip
1-1077	4-F-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
1-1078	4-F-Ph	2-MeNH-4-Pym	3-Hpip

Compound No.	R ¹	R ²	R ³
1-1079	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
1-1080	4-F-Ph	2-MeNH-4-Pym	4-Hpip
1-1081	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
1-1082	4-F-Ph	2-MeNH-4-Pym	2-Mor
1-1083	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor
1-1084	4-F-Ph	2-MeNH-4-Pym	2-Tmor
1-1085	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
1-1086	4-F-Ph	2-MeNH-4-Pym	1-Piz
1-1087	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1088	4-F-Ph	2-MeNH-4-Pym	2-Piz
1-1089	4-F-Ph	2-MeNH-4-Pym	4-Pyr
1-1090	4-F-Ph	2-MeNH-4-Pym	3-Pyr
1-1091	4-F-Ph	2-MeNH-4-Pym	2-Pyr
1-1092	4-F-Ph	2-MeNH-4-Pym	4-Pym
1-1093	4-F-Ph	2-MeNH-4-Pym	5-Pym
1-1094	4-F-Ph	2-MeNH-4-Pym	2-Pym
1-1095	4-F-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
1-1096	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
1-1097	4-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1098	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1099	4-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1100	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1101	4-F-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
1-1102	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
1-1103	4-F-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
1-1104	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
1-1105	4-F-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
1-1106	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
1-1107	4-F-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
1-1108	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
1-1109	4-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1110	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
1-1111	4-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1112	4-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1113	4-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1114	4-F-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
1-1115	4-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1116	4-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1117	4-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1118	3-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
1-1119	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
1-1120	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1121	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
1-1122	3-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂
1-1123	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
1-1124	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1125	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
1-1126	3-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂
1-1127	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
1-1128	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1129	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
1-1130	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
1-1131	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
1-1132	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1133	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
1-1134	3-F-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
1-1135	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
1-1136	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1137	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
1-1138	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-1139	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
1-1140	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1141	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
1-1142	3-F-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
1-1143	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
1-1144	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1145	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
1-1146	3-F-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
1-1147	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
1-1148	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1149	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
1-1150	3-F-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
1-1151	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
1-1152	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1153	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
1-1154	3-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1155	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
1-1156	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1157	3-F-Ph	2-MeNH-4-Pym	4-Mc-1-Piz-(CH ₂) ₃
1-1158	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
1-1159	3-F-Ph	2-MeNH-4-Pym	3-Azt
1-1160	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1161	3-F-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
1-1162	3-F-Ph	2-MeNH-4-Pym	3-Pyrd
1-1163	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1164	3-F-Ph	2-MeNH-4-Pym	3-Pip
1-1165	3-F-Ph	2-MeNH-4-Pym	4-Pip
1-1166	3-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1167	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1168	3-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-1169	3-F-Ph	2-MeNH-4-Pym	1-Et-4-Pip
1-1170	3-F-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
1-1171	3-F-Ph	2-MeNH-4-Pym	3-Hpip
1-1172	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
1-1173	3-F-Ph	2-MeNH-4-Pym	4-Hpip
1-1174	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
1-1175	3-F-Ph	2-MeNH-4-Pym	2-Mor
1-1176	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor
1-1177	3-F-Ph	2-MeNH-4-Pym	2-Tmor
1-1178	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
1-1179	3-F-Ph	2-MeNH-4-Pym	1-Piz
1-1180	3-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1181	3-F-Ph	2-MeNH-4-Pym	2-Piz
1-1182	3-F-Ph	2-MeNH-4-Pym	4-Pyr
1-1183	3-F-Ph	2-MeNH-4-Pym	3-Pyr
1-1184	3-F-Ph	2-MeNH-4-Pym	2-Pyr
1-1185	3-F-Ph	2-MeNH-4-Pym	4-Pym
1-1186	3-F-Ph	2-MeNH-4-Pym	5-Pym
1-1187	3-F-Ph	2-MeNH-4-Pym	2-Pym
1-1188	3-F-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
1-1189	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
1-1190	3-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1191	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1192	3-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1193	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1194	3-F-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
1-1195	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
1-1196	3-F-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
1-1197	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
1-1198	3-F-Ph	2-MeNH-4-Pym	2-Mor-CH ₂

Compound No.	R ¹	R ²	R ³
1-1199	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
1-1200	3-F-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
1-1201	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
1-1202	3-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1203	3-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
1-1204	3-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1205	3-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1206	3-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1207	3-F-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
1-1208	3-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1209	3-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1210	3-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1211	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
1-1212	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
1-1213	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1214	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
1-1215	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-CH ₂
1-1216	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
1-1217	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1218	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
1-1219	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-CH ₂
1-1220	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
1-1221	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1222	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
1-1223	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
1-1224	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
1-1225	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1226	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
1-1227	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
1-1228	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
1-1229	3,4-diF-Ph	2-McNH-4-Pym	1-Azt-(CH ₂) ₃
1-1230	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
1-1231	3,4-diF-Ph	2-McNH-4-Pym	1-Pyrd-CH ₂
1-1232	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
1-1233	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1234	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
1-1235	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
1-1236	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
1-1237	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1238	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
1-1239	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
1-1240	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
1-1241	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1242	3,4-diF-Ph	2-McNH-4-Pym	1-Mor-(CH ₂) ₄
1-1243	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
1-1244	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
1-1245	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1246	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
1-1247	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1248	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
1-1249	3,4-diF-Ph	2-McNH-4-Pym	1-Piz-(CH ₂) ₃
1-1250	3,4-diF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1251	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
1-1252	3,4-diF-Ph	2-MeNH-4-Pym	3-Azt
1-1253	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1254	3,4-diF-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
1-1255	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyrd
1-1256	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1257	3,4-diF-Ph	2-MeNH-4-Pym	3-Pip
1-1258	3,4-diF-Ph	2-MeNH-4-Pym	4-Pip

Compound No.	R ¹	R ²	R ³
1-1259	3,4-diF-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1260	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1261	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1262	3,4-diF-Ph	2-MeNH-4-Pym	1-Et-4-Pip
1-1263	3,4-diF-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
1-1264	3,4-diF-Ph	2-MeNH-4-Pym	3-Hpip
1-1265	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
1-1266	3,4-diF-Ph	2-MeNH-4-Pym	4-Hpip
1-1267	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
1-1268	3,4-diF-Ph	2-MeNH-4-Pym	2-Mor
1-1269	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Mor
1-1270	3,4-diF-Ph	2-MeNH-4-Pym	2-Tmor
1-1271	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
1-1272	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz
1-1273	3,4-diF-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1274	3,4-diF-Ph	2-MeNH-4-Pym	2-Piz
1-1275	3,4-diF-Ph	2-MeNH-4-Pym	4-Pyr
1-1276	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyr
1-1277	3,4-diF-Ph	2-MeNH-4-Pym	2-Pyr
1-1278	3,4-diF-Ph	2-MeNH-4-Pym	4-Pym
1-1279	3,4-diF-Ph	2-MeNH-4-Pym	5-Pym
1-1280	3,4-diF-Ph	2-MeNH-4-Pym	2-Pym
1-1281	3,4-diF-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
1-1282	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
1-1283	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1284	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1285	3,4-diF-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1286	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1287	3,4-diF-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
1-1288	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂

Compound No.	R ¹	R ²	R ³
1-1289	3,4-diF-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
1-1290	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
1-1291	3,4-diF-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
1-1292	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
1-1293	3,4-diF-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
1-1294	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
1-1295	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1296	3,4-diF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
1-1297	3,4-diF-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1298	3,4-diF-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1299	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1300	3,4-diF-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
1-1301	3,4-diF-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1302	3,4-diF-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1303	3,4-diF-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1304	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
1-1305	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
1-1306	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1307	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
1-1308	3-Cl-Ph	2-MeNH-4-Pym	MeNH-CH ₂
1-1309	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
1-1310	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1311	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
1-1312	3-Cl-Ph	2-MeNH-4-Pym	EtNH-CH ₂
1-1313	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
1-1314	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1315	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
1-1316	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
1-1317	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
1-1318	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1319	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
1-1320	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
1-1321	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
1-1322	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1323	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
1-1324	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
1-1325	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
1-1326	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1327	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
1-1328	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
1-1329	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
1-1330	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1331	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
1-1332	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
1-1333	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
1-1334	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1335	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
1-1336	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
1-1337	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
1-1338	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1339	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
1-1340	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1341	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
1-1342	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1343	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1344	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
1-1345	3-Cl-Ph	2-MeNH-4-Pym	3-Azt
1-1346	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1347	3-Cl-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
1-1348	3-Cl-Ph	2-MeNH-4-Pym	3-Pyrd

Compound No.	R ¹	R ²	R ³
1-1349	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1350	3-Cl-Ph	2-MeNH-4-Pym	3-Pip
1-1351	3-Cl-Ph	2-MeNH-4-Pym	4-Pip
1-1352	3-Cl-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1353	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1354	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1355	3-Cl-Ph	2-MeNH-4-Pym	1-Et-4-Pip
1-1356	3-Cl-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
1-1357	3-Cl-Ph	2-MeNH-4-Pym	3-Hpip
1-1358	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
1-1359	3-Cl-Ph	2-MeNH-4-Pym	4-Hpip
1-1360	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
1-1361	3-Cl-Ph	2-MeNH-4-Pym	2-Mor
1-1362	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Mor
1-1363	3-Cl-Ph	2-MeNH-4-Pym	2-Tmor
1-1364	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
1-1365	3-Cl-Ph	2-MeNH-4-Pym	1-Piz
1-1366	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1367	3-Cl-Ph	2-MeNH-4-Pym	2-Piz
1-1368	3-Cl-Ph	2-MeNH-4-Pym	4-Pyr
1-1369	3-Cl-Ph	2-MeNH-4-Pym	3-Pyr
1-1370	3-Cl-Ph	2-MeNH-4-Pym	2-Pyr
1-1371	3-Cl-Ph	2-MeNH-4-Pym	4-Pym
1-1372	3-Cl-Ph	2-MeNH-4-Pym	5-Pym
1-1373	3-Cl-Ph	2-MeNH-4-Pym	2-Pym
1-1374	3-Cl-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
1-1375	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
1-1376	3-Cl-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1377	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1378	3-Cl-Ph	2-MeNH-4-Pym	4-Pip-CH ₂

Compound No.	R ¹	R ²	R ³
1-1379	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1380	3-Cl-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
1-1381	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
1-1382	3-Cl-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
1-1383	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
1-1384	3-Cl-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
1-1385	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
1-1386	3-Cl-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
1-1387	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
1-1388	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
1-1389	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
1-1390	3-Cl-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1391	3-Cl-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1392	3-Cl-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1393	3-Cl-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
1-1394	3-Cl-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1395	3-Cl-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1396	3-Cl-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1397	3-Cl-4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-1398	3-Cl-4-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-1399	3-Cl-4-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-1400	3-Cl-4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-1401	3-Cl-4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-1402	3-Cl-4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-1403	3-Cl-4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-1404	3-Cl-4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-1405	3-Cl-4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-1406	3-Cl-4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-1407	3-Cl-4-F-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1408	3-Cl-4-F-Ph	4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
1-1409	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Azt
1-1410	3-Cl-4-F-Ph	4-Pyr	3-Pyrd
1-1411	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Pyrd
1-1412	3-Cl-4-F-Ph	4-Pyr	4-Pip
1-1413	3-Cl-4-F-Ph	4-Pyr	4-(3,4-deH-Pip)
1-1414	3-Cl-4-F-Ph	4-Pyr	1-Me-4-Pip
1-1415	3-Cl-4-F-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1416	3-Cl-4-F-Ph	4-Pyr	1-Piz
1-1417	3-Cl-4-F-Ph	4-Pyr	4-Me-1-Piz
1-1418	3-Cl-4-F-Ph	4-Pyr	4-Pyr
1-1419	3-Cl-4-F-Ph	4-Pyr	3-Pyr
1-1420	3-Cl-4-F-Ph	4-Pyr	4-Pym
1-1421	3-Cl-4-F-Ph	4-Pyr	5-Pym
1-1422	3-Cl-4-F-Ph	4-Pyr	3-Pyrd-CH ₂
1-1423	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-1424	3-Cl-4-F-Ph	4-Pyr	4-Pip-CH ₂
1-1425	3-Cl-4-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-1426	3-Cl-4-F-Ph	4-Pyr	2-Piz-CH ₂
1-1427	3-Cl-4-F-Ph	4-Pyr	4-Pyr-CH ₂
1-1428	3-Cl-4-F-Ph	4-Pyr	3-Pyr-CH ₂
1-1429	3-Cl-4-F-Ph	4-Pyr	4-Pym-CH ₂
1-1430	3-Cl-4-F-Ph	4-Pyr	5-Pym-CH ₂
1-1431	3-Cl-4-F-Ph	4-Pyr	2-Pym-CH ₂
1-1432	3,4,5-triF-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-1433	3,4,5-triF-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-1434	3,4,5-triF-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-1435	3,4,5-triF-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-1436	3,4,5-triF-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-1437	3,4,5-triF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-1438	3,4,5-triF-Ph	4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1439	3,4,5-triF-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-1440	3,4,5-triF-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-1441	3,4,5-triF-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-1442	3,4,5-triF-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1443	3,4,5-triF-Ph	4-Pyr	3-Azt
1-1444	3,4,5-triF-Ph	4-Pyr	1-Me-3-Azt
1-1445	3,4,5-triF-Ph	4-Pyr	3-Pyrd
1-1446	3,4,5-triF-Ph	4-Pyr	1-Me-3-Pyrd
1-1447	3,4,5-triF-Ph	4-Pyr	4-Pip
1-1448	3,4,5-triF-Ph	4-Pyr	4-(3,4-deH-Pip)
1-1449	3,4,5-triF-Ph	4-Pyr	1-Me-4-Pip
1-1450	3,4,5-triF-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1451	3,4,5-triF-Ph	4-Pyr	1-Piz
1-1452	3,4,5-triF-Ph	4-Pyr	4-Me-1-Piz
1-1453	3,4,5-triF-Ph	4-Pyr	4-Pyr
1-1454	3,4,5-triF-Ph	4-Pyr	3-Pyr
1-1455	3,4,5-triF-Ph	4-Pyr	4-Pym
1-1456	3,4,5-triF-Ph	4-Pyr	5-Pym
1-1457	3,4,5-triF-Ph	4-Pyr	3-Pyrd-CH ₂
1-1458	3,4,5-triF-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-1459	3,4,5-triF-Ph	4-Pyr	4-Pip-CH ₂
1-1460	3,4,5-triF-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-1461	3,4,5-triF-Ph	4-Pyr	2-Piz-CH ₂
1-1462	3,4,5-triF-Ph	4-Pyr	4-Pyr-CH ₂
1-1463	3,4,5-triF-Ph	4-Pyr	3-Pyr-CH ₂
1-1464	3,4,5-triF-Ph	4-Pyr	4-Pym-CH ₂
1-1465	3,4,5-triF-Ph	4-Pyr	5-Pym-CH ₂
1-1466	3,4,5-triF-Ph	4-Pyr	2-Pym-CH ₂
1-1467	3-CF ₃ -Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-1468	3-CF ₃ -Ph	4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1469	3-CF ₃ -Ph	4-Pyr	EtNH-(CH ₂) ₃
1-1470	3-CF ₃ -Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-1471	3-CF ₃ -Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-1472	3-CF ₃ -Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-1473	3-CF ₃ -Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-1474	3-CF ₃ -Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-1475	3-CF ₃ -Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-1476	3-CF ₃ -Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-1477	3-CF ₃ -Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1478	3-CF ₃ -Ph	4-Pyr	3-Azt
1-1479	3-CF ₃ -Ph	4-Pyr	1-Me-3-Azt
1-1480	3-CF ₃ -Ph	4-Pyr	3-Pyrd
1-1481	3-CF ₃ -Ph	4-Pyr	1-Me-3-Pyrd
1-1482	3-CF ₃ -Ph	4-Pyr	4-Pip
1-1483	3-CF ₃ -Ph	4-Pyr	4-(3,4-deH-Pip)
1-1484	3-CF ₃ -Ph	4-Pyr	1-Me-4-Pip
1-1485	3-CF ₃ -Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1486	3-CF ₃ -Ph	4-Pyr	1-Piz
1-1487	3-CF ₃ -Ph	4-Pyr	4-Me-1-Piz
1-1488	3-CF ₃ -Ph	4-Pyr	4-Pyr
1-1489	3-CF ₃ -Ph	4-Pyr	3-Pyr
1-1490	3-CF ₃ -Ph	4-Pyr	4-Pym
1-1491	3-CF ₃ -Ph	4-Pyr	5-Pym
1-1492	3-CF ₃ -Ph	4-Pyr	3-Pyrd-CH ₂
1-1493	3-CF ₃ -Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
1-1494	3-CF ₃ -Ph	4-Pyr	4-Pip-CH ₂
1-1495	3-CF ₃ -Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-1496	3-CF ₃ -Ph	4-Pyr	2-Piz-CH ₂
1-1497	3-CF ₃ -Ph	4-Pyr	4-Pyr-CH ₂
1-1498	3-CF ₃ -Ph	4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-1499	3-CF ₃ -Ph	4-Pyr	4-Pym-CH ₂
1-1500	3-CF ₃ -Ph	4-Pyr	5-Pym-CH ₂
1-1501	3-CF ₃ -Ph	4-Pyr	2-Pym-CH ₂
1-1502	3-CHF ₂ O-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
1-1503	3-CHF ₂ O-Ph	4-Pyr	MeNH-(CH ₂) ₃
1-1504	3-CHF ₂ O-Ph	4-Pyr	EtNH-(CH ₂) ₃
1-1505	3-CHF ₂ O-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
1-1506	3-CHF ₂ O-Ph	4-Pyr	1-Azt-(CH ₂) ₃
1-1507	3-CHF ₂ O-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
1-1508	3-CHF ₂ O-Ph	4-Pyr	1-Pip-(CH ₂) ₃
1-1509	3-CHF ₂ O-Ph	4-Pyr	1-Mor-(CH ₂) ₃
1-1510	3-CHF ₂ O-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
1-1511	3-CHF ₂ O-Ph	4-Pyr	1-Piz-(CH ₂) ₃
1-1512	3-CHF ₂ O-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1513	3-CHF ₂ O-Ph	4-Pyr	3-Azt
1-1514	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Azt
1-1515	3-CHF ₂ O-Ph	4-Pyr	3-Pyrd
1-1516	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Pyrd
1-1517	3-CHF ₂ O-Ph	4-Pyr	4-Pip
1-1518	3-CHF ₂ O-Ph	4-Pyr	4-(3,4-deH-Pip)
1-1519	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-Pip
1-1520	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1521	3-CHF ₂ O-Ph	4-Pyr	1-Piz
1-1522	3-CHF ₂ O-Ph	4-Pyr	4-Me-1-Piz
1-1523	3-CHF ₂ O-Ph	4-Pyr	4-Pyr
1-1524	3-CHF ₂ O-Ph	4-Pyr	3-Pyr
1-1525	3-CHF ₂ O-Ph	4-Pyr	4-Pym
1-1526	3-CHF ₂ O-Ph	4-Pyr	5-Pym
1-1527	3-CHF ₂ O-Ph	4-Pyr	3-Pyrd-CH ₂
1-1528	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-1529	3-CHF ₂ O-Ph	4-Pyr	4-Pip-CH ₂
1-1530	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-Pip-CH ₂
1-1531	3-CHF ₂ O-Ph	4-Pyr	2-Piz-CH ₂
1-1532	3-CHF ₂ O-Ph	4-Pyr	4-Pyr-CH ₂
1-1533	3-CHF ₂ O-Ph	4-Pyr	3-Pyr-CH ₂
1-1534	3-CHF ₂ O-Ph	4-Pyr	4-Pym-CH ₂
1-1535	3-CHF ₂ O-Ph	4-Pyr	5-Pym-CH ₂
1-1536	3-CHF ₂ O-Ph	4-Pyr	2-Pym-CH ₂
1-1537	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-1538	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-1539	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-1540	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-1541	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-1542	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-1543	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-1544	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-1545	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-1546	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-1547	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1548	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Azt
1-1549	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-1550	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-1551	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-1552	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pip
1-1553	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-1554	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-1555	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-1556	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Piz
1-1557	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-1558	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
1-1559	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyr
1-1560	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pym
1-1561	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	5-Pym
1-1562	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-1563	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-1564	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-1565	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-1566	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-1567	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-1568	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-1569	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-1570	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-1571	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-1572	3,4,5-triF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-1573	3,4,5-triF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-1574	3,4,5-triF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-1575	3,4,5-triF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-1576	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-1577	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-1578	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-1579	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-1580	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-1581	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-1582	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1583	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Azt
1-1584	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-1585	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-1586	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-1587	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pip
1-1588	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-1589	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-1590	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-d ^e H-Pip)
1-1591	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Piz
1-1592	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-1593	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pyr
1-1594	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyr
1-1595	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pym
1-1596	3,4,5-triF-Ph	2-NH ₂ -4-Pym	5-Pym
1-1597	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-1598	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-1599	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-1600	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-1601	3,4,5-triF-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-1602	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-1603	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-1604	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-1605	3,4,5-triF-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-1606	3,4,5-triF-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-1607	3-CF ₃ -Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-1608	3-CF ₃ -Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-1609	3-CF ₃ -Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-1610	3-CF ₃ -Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-1611	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-1612	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-1613	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
1-1614	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-1615	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-1616	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-1617	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₂
1-1618	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
1-1619	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-1620	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyrd
1-1621	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-1622	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pip
1-1623	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-1624	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-1625	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-1626	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Piz
1-1627	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-1628	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pyr
1-1629	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyr
1-1630	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pym
1-1631	3-CF ₃ -Ph	2-NH ₂ -4-Pym	5-Pym
1-1632	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-1633	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-1634	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-1635	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-1636	3-CF ₃ -Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-1637	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-1638	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-1639	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-1640	3-CF ₃ -Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-1641	3-CF ₃ -Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-1642	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
1-1643	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
1-1644	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
1-1645	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
1-1646	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
1-1647	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
1-1648	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1649	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
1-1650	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
1-1651	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
1-1652	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1653	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Azt
1-1654	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
1-1655	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyrd
1-1656	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
1-1657	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pip
1-1658	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
1-1659	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
1-1660	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
1-1661	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Piz
1-1662	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
1-1663	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pyr
1-1664	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyr
1-1665	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pym
1-1666	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	5-Pym
1-1667	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
1-1668	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
1-1669	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
1-1670	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
1-1671	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
1-1672	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
1-1673	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
1-1674	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
1-1675	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
1-1676	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
1-1677	3-Cl-4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1678	3-Cl-4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1679	3-Cl-4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1680	3-Cl-4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1681	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1682	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1683	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1684	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1685	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1686	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1687	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1688	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Azt
1-1689	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1690	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyrd
1-1691	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1692	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pip
1-1693	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1694	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1695	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1696	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Piz
1-1697	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1698	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pyr
1-1699	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyr
1-1700	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pym
1-1701	3-Cl-4-F-Ph	2-MeNH-4-Pym	5-Pym
1-1702	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1703	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1704	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1705	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1706	3-Cl-4-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1707	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1708	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-1709	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1710	3-Cl-4-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1711	3-Cl-4-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1712	3,4,5-triF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1713	3,4,5-triF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1714	3,4,5-triF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1715	3,4,5-triF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1716	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1717	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1718	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1719	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1720	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1721	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1722	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1723	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Azt
1-1724	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1725	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyrd
1-1726	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1727	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pip
1-1728	3,4,5-triF-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1729	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1730	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1731	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Piz
1-1732	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1733	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pyr
1-1734	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyr
1-1735	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pym
1-1736	3,4,5-triF-Ph	2-MeNH-4-Pym	5-Pym
1-1737	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1738	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-1739	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1740	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1741	3,4,5-triF-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1742	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1743	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1744	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1745	3,4,5-triF-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1746	3,4,5-triF-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1747	3-CF ₃ -Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1748	3-CF ₃ -Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1749	3-CF ₃ -Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1750	3-CF ₃ -Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1751	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1752	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1753	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1754	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1755	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1756	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1757	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1758	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Azt
1-1759	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1760	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyrd
1-1761	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1762	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pip
1-1763	3-CF ₃ -Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
1-1764	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1765	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1766	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Piz
1-1767	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1768	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
1-1769	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyr
1-1770	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pym
1-1771	3-CF ₃ -Ph	2-MeNH-4-Pym	5-Pym
1-1772	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1773	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1774	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1775	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1776	3-CF ₃ -Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1777	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1778	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1779	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1780	3-CF ₃ -Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1781	3-CF ₃ -Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1782	3-CHF ₂ O-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
1-1783	3-CHF ₂ O-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
1-1784	3-CHF ₂ O-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
1-1785	3-CHF ₂ O-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
1-1786	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
1-1787	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
1-1788	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
1-1789	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
1-1790	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
1-1791	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
1-1792	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-1793	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Azt
1-1794	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Azt
1-1795	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyrd
1-1796	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
1-1797	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pip
1-1798	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-1799	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-Pip
1-1800	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
1-1801	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Piz
1-1802	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Me-1-Piz
1-1803	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pyr
1-1804	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyr
1-1805	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pym
1-1806	3-CHF ₂ O-Ph	2-MeNH-4-Pym	5-Pym
1-1807	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
1-1808	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
1-1809	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
1-1810	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
1-1811	3-CHF ₂ O-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
1-1812	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
1-1813	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
1-1814	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
1-1815	3-CHF ₂ O-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
1-1816	3-CHF ₂ O-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
1-1817	Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
1-1818	Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
1-1819	Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
1-1820	Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
1-1821	Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
1-1822	Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
1-1823	Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
1-1824	Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
1-1825	Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
1-1826	Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
1-1827	Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1828	Ph	2-NH ₂ -4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
1-1829	Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
1-1830	Ph	2-NH ₂ -4-Pyr	3-Pyrd
1-1831	Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
1-1832	Ph	2-NH ₂ -4-Pyr	4-Pip
1-1833	Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
1-1834	Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
1-1835	Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1836	Ph	2-NH ₂ -4-Pyr	1-Piz
1-1837	Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
1-1838	Ph	2-NH ₂ -4-Pyr	4-Pyr
1-1839	Ph	2-NH ₂ -4-Pyr	3-Pyr
1-1840	Ph	2-NH ₂ -4-Pyr	4-Pym
1-1841	Ph	2-NH ₂ -4-Pyr	5-Pym
1-1842	Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
1-1843	Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
1-1844	Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
1-1845	Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
1-1846	Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
1-1847	Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
1-1848	Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
1-1849	Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
1-1850	Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
1-1851	Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
1-1852	4-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
1-1853	4-F-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
1-1854	4-F-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
1-1855	4-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
1-1856	4-F-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
1-1857	4-F-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
1-1858	4-F-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1859	4-F-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
1-1860	4-F-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
1-1861	4-F-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
1-1862	4-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1863	4-F-Ph	2-NH ₂ -4-Pyr	3-Azt
1-1864	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
1-1865	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd
1-1866	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
1-1867	4-F-Ph	2-NH ₂ -4-Pyr	4-Pip
1-1868	4-F-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
1-1869	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
1-1870	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1871	4-F-Ph	2-NH ₂ -4-Pyr	1-Piz
1-1872	4-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
1-1873	4-F-Ph	2-NH ₂ -4-Pyr	4-Pyr
1-1874	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyr
1-1875	4-F-Ph	2-NH ₂ -4-Pyr	4-Pym
1-1876	4-F-Ph	2-NH ₂ -4-Pyr	5-Pym
1-1877	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
1-1878	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
1-1879	4-F-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
1-1880	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
1-1881	4-F-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
1-1882	4-F-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
1-1883	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
1-1884	4-F-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
1-1885	4-F-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
1-1886	4-F-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
1-1887	3-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
1-1888	3-F-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-1889	3-F-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
1-1890	3-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
1-1891	3-F-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
1-1892	3-F-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
1-1893	3-F-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
1-1894	3-F-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
1-1895	3-F-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
1-1896	3-F-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
1-1897	3-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1898	3-F-Ph	2-NH ₂ -4-Pyr	3-Azt
1-1899	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
1-1900	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd
1-1901	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
1-1902	3-F-Ph	2-NH ₂ -4-Pyr	4-Pip
1-1903	3-F-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
1-1904	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
1-1905	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1906	3-F-Ph	2-NH ₂ -4-Pyr	1-Piz
1-1907	3-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
1-1908	3-F-Ph	2-NH ₂ -4-Pyr	4-Pyr
1-1909	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyr
1-1910	3-F-Ph	2-NH ₂ -4-Pyr	4-Pym
1-1911	3-F-Ph	2-NH ₂ -4-Pyr	5-Pym
1-1912	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
1-1913	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
1-1914	3-F-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
1-1915	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
1-1916	3-F-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
1-1917	3-F-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
1-1918	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-1919	3-F-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
1-1920	3-F-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
1-1921	3-F-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
1-1922	3,4-diF-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
1-1923	3,4-diF-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
1-1924	3,4-diF-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
1-1925	3,4-diF-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
1-1926	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
1-1927	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
1-1928	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
1-1929	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
1-1930	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
1-1931	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
1-1932	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1933	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Azt
1-1934	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
1-1935	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyrd
1-1936	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
1-1937	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pip
1-1938	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
1-1939	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
1-1940	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1941	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Piz
1-1942	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
1-1943	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pyr
1-1944	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyr
1-1945	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pym
1-1946	3,4-diF-Ph	2-NH ₂ -4-Pyr	5-Pym
1-1947	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
1-1948	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-1949	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
1-1950	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
1-1951	3,4-diF-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
1-1952	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
1-1953	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
1-1954	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
1-1955	3,4-diF-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
1-1956	3,4-diF-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
1-1957	3-Cl-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
1-1958	3-Cl-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
1-1959	3-Cl-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
1-1960	3-Cl-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
1-1961	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
1-1962	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
1-1963	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
1-1964	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
1-1965	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
1-1966	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
1-1967	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-1968	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Azt
1-1969	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
1-1970	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyrd
1-1971	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
1-1972	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pip
1-1973	3-Cl-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
1-1974	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
1-1975	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
1-1976	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Piz
1-1977	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
1-1978	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pyr

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Compound No.	R ¹	R ²	R ³
1-1979	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyr
1-1980	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pym
1-1981	3-Cl-Ph	2-NH ₂ -4-Pyr	5-Pym
1-1982	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
1-1983	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
1-1984	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
1-1985	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
1-1986	3-Cl-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
1-1987	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
1-1988	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
1-1989	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
1-1990	3-Cl-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
1-1991	3-Cl-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
1-1992	Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
1-1993	Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
1-1994	Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
1-1995	Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
1-1996	Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
1-1997	Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
1-1998	Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
1-1999	Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
1-2000	Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
1-2001	Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
1-2002	Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-2003	Ph	2-MeNH-4-Pyr	3-Azt
1-2004	Ph	2-MeNH-4-Pyr	1-Me-3-Azt
1-2005	Ph	2-MeNH-4-Pyr	3-Pyrd
1-2006	Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
1-2007	Ph	2-MeNH-4-Pyr	4-Pip
1-2008	Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2009	Ph	2-MeNH-4-Pyr	1-Me-4-Pip
1-2010	Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
1-2011	Ph	2-MeNH-4-Pyr	1-Piz
1-2012	Ph	2-MeNH-4-Pyr	4-Me-1-Piz
1-2013	Ph	2-MeNH-4-Pyr	4-Pyr
1-2014	Ph	2-MeNH-4-Pyr	3-Pyr
1-2015	Ph	2-MeNH-4-Pyr	4-Pym
1-2016	Ph	2-MeNH-4-Pyr	5-Pym
1-2017	Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
1-2018	Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
1-2019	Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
1-2020	Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
1-2021	Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
1-2022	Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
1-2023	Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
1-2024	Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
1-2025	Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
1-2026	Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
1-2027	4-F-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
1-2028	4-F-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
1-2029	4-F-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
1-2030	4-F-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
1-2031	4-F-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
1-2032	4-F-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
1-2033	4-F-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
1-2034	4-F-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
1-2035	4-F-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
1-2036	4-F-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
1-2037	4-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-2038	4-F-Ph	2-MeNH-4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
1-2039	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
1-2040	4-F-Ph	2-MeNH-4-Pyr	3-Pyrd
1-2041	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
1-2042	4-F-Ph	2-MeNH-4-Pyr	4-Pip
1-2043	4-F-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
1-2044	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
1-2045	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
1-2046	4-F-Ph	2-MeNH-4-Pyr	1-Piz
1-2047	4-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
1-2048	4-F-Ph	2-MeNH-4-Pyr	4-Pyr
1-2049	4-F-Ph	2-MeNH-4-Pyr	3-Pyr
1-2050	4-F-Ph	2-MeNH-4-Pyr	4-Pym
1-2051	4-F-Ph	2-MeNH-4-Pyr	5-Pym
1-2052	4-F-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
1-2053	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
1-2054	4-F-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
1-2055	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
1-2056	4-F-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
1-2057	4-F-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
1-2058	4-F-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
1-2059	4-F-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
1-2060	4-F-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
1-2061	4-F-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
1-2062	3-F-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
1-2063	3-F-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
1-2064	3-F-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
1-2065	3-F-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
1-2066	3-F-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
1-2067	3-F-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
1-2068	3-F-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-2069	3-F-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
1-2070	3-F-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
1-2071	3-F-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
1-2072	3-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-2073	3-F-Ph	2-MeNH-4-Pyr	3-Azt
1-2074	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
1-2075	3-F-Ph	2-MeNH-4-Pyr	3-Pyrd
1-2076	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
1-2077	3-F-Ph	2-MeNH-4-Pyr	4-Pip
1-2078	3-F-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
1-2079	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
1-2080	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
1-2081	3-F-Ph	2-MeNH-4-Pyr	1-Piz
1-2082	3-F-Ph	2-MeNH-4-Pyr	4-Mc-1-Piz
1-2083	3-F-Ph	2-MeNH-4-Pyr	4-Pyr
1-2084	3-F-Ph	2-MeNH-4-Pyr	3-Pyr
1-2085	3-F-Ph	2-MeNH-4-Pyr	4-Pym
1-2086	3-F-Ph	2-MeNH-4-Pyr	5-Pym
1-2087	3-F-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
1-2088	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
1-2089	3-F-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
1-2090	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
1-2091	3-F-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
1-2092	3-F-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
1-2093	3-F-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
1-2094	3-F-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
1-2095	3-F-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
1-2096	3-F-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
1-2097	3,4-diF-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
1-2098	3,4-diF-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-2099	3,4-diF-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
1-2100	3,4-diF-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
1-2101	3,4-diF-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
1-2102	3,4-diF-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
1-2103	3,4-diF-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
1-2104	3,4-diF-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
1-2105	3,4-diF-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
1-2106	3,4-diF-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
1-2107	3,4-diF-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-2108	3,4-diF-Ph	2-MeNH-4-Pyr	3-Azt
1-2109	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
1-2110	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyrd
1-2111	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
1-2112	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pip
1-2113	3,4-diF-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
1-2114	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
1-2115	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
1-2116	3,4-diF-Ph	2-MeNH-4-Pyr	1-Piz
1-2117	3,4-diF-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
1-2118	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pyr
1-2119	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyr
1-2120	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pym
1-2121	3,4-diF-Ph	2-MeNH-4-Pyr	5-Pym
1-2122	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
1-2123	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
1-2124	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
1-2125	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
1-2126	3,4-diF-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
1-2127	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
1-2128	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-2129	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
1-2130	3,4-diF-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
1-2131	3,4-diF-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
1-2132	3-Cl-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
1-2133	3-Cl-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
1-2134	3-Cl-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
1-2135	3-Cl-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
1-2136	3-Cl-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
1-2137	3-Cl-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
1-2138	3-Cl-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
1-2139	3-Cl-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
1-2140	3-Cl-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
1-2141	3-Cl-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
1-2142	3-Cl-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
1-2143	3-Cl-Ph	2-MeNH-4-Pyr	3-Azt
1-2144	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
1-2145	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyrd
1-2146	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
1-2147	3-Cl-Ph	2-MeNH-4-Pyr	4-Pip
1-2148	3-Cl-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
1-2149	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
1-2150	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
1-2151	3-Cl-Ph	2-MeNH-4-Pyr	1-Piz
1-2152	3-Cl-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
1-2153	3-Cl-Ph	2-MeNH-4-Pyr	4-Pyr
1-2154	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyr
1-2155	3-Cl-Ph	2-MeNH-4-Pyr	4-Pym
1-2156	3-Cl-Ph	2-MeNH-4-Pyr	5-Pym
1-2157	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
1-2158	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-2159	3-Cl-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
1-2160	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
1-2161	3-Cl-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
1-2162	3-Cl-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
1-2163	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
1-2164	3-Cl-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
1-2165	3-Cl-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
1-2166	3-Cl-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
1-2167	Ph	4-Pym	H ₂ N-(CH ₂) ₃
1-2168	Ph	4-Pym	MeNH-(CH ₂) ₃
1-2169	Ph	4-Pym	EtNH-(CH ₂) ₃
1-2170	Ph	4-Pym	Me ₂ N-(CH ₂) ₃
1-2171	Ph	4-Pym	1-Azt-(CH ₂) ₃
1-2172	Ph	4-Pym	1-Pyrd-(CH ₂) ₃
1-2173	Ph	4-Pym	1-Pip-(CH ₂) ₃
1-2174	Ph	4-Pym	1-Mor-(CH ₂) ₃
1-2175	Ph	4-Pym	1-Tmor-(CH ₂) ₃
1-2176	Ph	4-Pym	1-Piz-(CH ₂) ₃
1-2177	Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2178	Ph	4-Pym	3-Azt
1-2179	Ph	4-Pym	1-Me-3-Azt
1-2180	Ph	4-Pym	3-Pyrd
1-2181	Ph	4-Pym	1-Me-3-Pyrd
1-2182	Ph	4-Pym	4-Pip
1-2183	Ph	4-Pym	4-(3,4-deH-Pip)
1-2184	Ph	4-Pym	1-Me-4-Pip
1-2185	Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
1-2186	Ph	4-Pym	1-Piz
1-2187	Ph	4-Pym	4-Me-1-Piz
1-2188	Ph	4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
1-2189	Ph	4-Pym	3-Pyr
1-2190	Ph	4-Pym	4-Pym
1-2191	Ph	4-Pym	5-Pym
1-2192	Ph	4-Pym	3-Pyrd-CH ₂
1-2193	Ph	4-Pym	1-Me-3-Pyrd-CH ₂
1-2194	Ph	4-Pym	4-Pip-CH ₂
1-2195	Ph	4-Pym	1-Me-4-Pip-CH ₂
1-2196	Ph	4-Pym	2-Piz-CH ₂
1-2197	Ph	4-Pym	4-Pyr-CH ₂
1-2198	Ph	4-Pym	3-Pyr-CH ₂
1-2199	Ph	4-Pym	4-Pym-CH ₂
1-2200	Ph	4-Pym	5-Pym-CH ₂
1-2201	Ph	4-Pym	2-Pym-CH ₂
1-2202	4-F-Ph	4-Pym	H ₂ N-(CH ₂) ₃
1-2203	4-F-Ph	4-Pym	MeNH-(CH ₂) ₃
1-2204	4-F-Ph	4-Pym	EtNH-(CH ₂) ₃
1-2205	4-F-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
1-2206	4-F-Ph	4-Pym	1-Azt-(CH ₂) ₃
1-2207	4-F-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
1-2208	4-F-Ph	4-Pym	1-Pip-(CH ₂) ₃
1-2209	4-F-Ph	4-Pym	1-Mor-(CH ₂) ₃
1-2210	4-F-Ph	4-Pym	1-Tmor-(CH ₂) ₃
1-2211	4-F-Ph	4-Pym	1-Piz-(CH ₂) ₃
1-2212	4-F-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2213	4-F-Ph	4-Pym	3-Azt
1-2214	4-F-Ph	4-Pym	1-Me-3-Azt
1-2215	4-F-Ph	4-Pym	3-Pyrd
1-2216	4-F-Ph	4-Pym	1-Me-3-Pyrd
1-2217	4-F-Ph	4-Pym	4-Pip
1-2218	4-F-Ph	4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2219	4-F-Ph	4-Pym	1-Me-4-Pip
1-2220	4-F-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
1-2221	4-F-Ph	4-Pym	1-Piz
1-2222	4-F-Ph	4-Pym	4-Me-1-Piz
1-2223	4-F-Ph	4-Pym	4-Pyr
1-2224	4-F-Ph	4-Pym	3-Pyr
1-2225	4-F-Ph	4-Pym	4-Pym
1-2226	4-F-Ph	4-Pym	5-Pym
1-2227	4-F-Ph	4-Pym	3-Pyrd-CH ₂
1-2228	4-F-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
1-2229	4-F-Ph	4-Pym	4-Pip-CH ₂
1-2230	4-F-Ph	4-Pym	1-Me-4-Pip-CH ₂
1-2231	4-F-Ph	4-Pym	2-Piz-CH ₂
1-2232	4-F-Ph	4-Pym	4-Pyr-CH ₂
1-2233	4-F-Ph	4-Pym	3-Pyr-CH ₂
1-2234	4-F-Ph	4-Pym	4-Pym-CH ₂
1-2235	4-F-Ph	4-Pym	5-Pym-CH ₂
1-2236	4-F-Ph	4-Pym	2-Pym-CH ₂
1-2237	3-F-Ph	4-Pym	H ₂ N-(CH ₂) ₃
1-2238	3-F-Ph	4-Pym	MeNH-(CH ₂) ₃
1-2239	3-F-Ph	4-Pym	EtNH-(CH ₂) ₃
1-2240	3-F-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
1-2241	3-F-Ph	4-Pym	1-Azt-(CH ₂) ₃
1-2242	3-F-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
1-2243	3-F-Ph	4-Pym	1-Pip-(CH ₂) ₃
1-2244	3-F-Ph	4-Pym	1-Mor-(CH ₂) ₃
1-2245	3-F-Ph	4-Pym	1-Tmor-(CH ₂) ₃
1-2246	3-F-Ph	4-Pym	1-Piz-(CH ₂) ₃
1-2247	3-F-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2248	3-F-Ph	4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
1-2249	3-F-Ph	4-Pym	1-Me-3-Azt
1-2250	3-F-Ph	4-Pym	3-Pyrd
1-2251	3-F-Ph	4-Pym	1-Me-3-Pyrd
1-2252	3-F-Ph	4-Pym	4-Pip
1-2253	3-F-Ph	4-Pym	4-(3,4-deH-Pip)
1-2254	3-F-Ph	4-Pym	1-Me-4-Pip
1-2255	3-F-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
1-2256	3-F-Ph	4-Pym	1-Piz
1-2257	3-F-Ph	4-Pym	4-Me-1-Piz
1-2258	3-F-Ph	4-Pym	4-Pyr
1-2259	3-F-Ph	4-Pym	3-Pyr
1-2260	3-F-Ph	4-Pym	4-Pym
1-2261	3-F-Ph	4-Pym	5-Pym
1-2262	3-F-Ph	4-Pym	3-Pyrd-CH ₂
1-2263	3-F-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
1-2264	3-F-Ph	4-Pym	4-Pip-CH ₂
1-2265	3-F-Ph	4-Pym	1-Me-4-Pip-CH ₂
1-2266	3-F-Ph	4-Pym	2-Piz-CH ₂
1-2267	3-F-Ph	4-Pym	4-Pyr-CH ₂
1-2268	3-F-Ph	4-Pym	3-Pyr-CH ₂
1-2269	3-F-Ph	4-Pym	4-Pym-CH ₂
1-2270	3-F-Ph	4-Pym	5-Pym-CH ₂
1-2271	3-F-Ph	4-Pym	2-Pym-CH ₂
1-2272	3,4-diF-Ph	4-Pym	H ₂ N-(CH ₂) ₃
1-2273	3,4-diF-Ph	4-Pym	MeNH-(CH ₂) ₃
1-2274	3,4-diF-Ph	4-Pym	EtNH-(CH ₂) ₃
1-2275	3,4-diF-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
1-2276	3,4-diF-Ph	4-Pym	1-Azt-(CH ₂) ₃
1-2277	3,4-diF-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
1-2278	3,4-diF-Ph	4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-2279	3,4-diF-Ph	4-Pym	1-Mor-(CH ₂) ₃
1-2280	3,4-diF-Ph	4-Pym	1-Tmor-(CH ₂) ₃
1-2281	3,4-diF-Ph	4-Pym	1-Piz-(CH ₂) ₃
1-2282	3,4-diF-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2283	3,4-diF-Ph	4-Pym	3-Azt
1-2284	3,4-diF-Ph	4-Pym	1-Me-3-Azt
1-2285	3,4-diF-Ph	4-Pym	3-Pyrd
1-2286	3,4-diF-Ph	4-Pym	1-Me-3-Pyrd
1-2287	3,4-diF-Ph	4-Pym	4-Pip
1-2288	3,4-diF-Ph	4-Pym	4-(3,4-deH-Pip)
1-2289	3,4-diF-Ph	4-Pym	1-Me-4-Pip
1-2290	3,4-diF-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
1-2291	3,4-diF-Ph	4-Pym	1-Piz
1-2292	3,4-diF-Ph	4-Pym	4-Me-1-Piz
1-2293	3,4-diF-Ph	4-Pym	4-Pyr
1-2294	3,4-diF-Ph	4-Pym	3-Pyr
1-2295	3,4-diF-Ph	4-Pym	4-Pym
1-2296	3,4-diF-Ph	4-Pym	5-Pym
1-2297	3,4-diF-Ph	4-Pym	3-Pyrd-CH ₂
1-2298	3,4-diF-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
1-2299	3,4-diF-Ph	4-Pym	4-Pip-CH ₂
1-2300	3,4-diF-Ph	4-Pym	1-Me-4-Pip-CH ₂
1-2301	3,4-diF-Ph	4-Pym	2-Piz-CH ₂
1-2302	3,4-diF-Ph	4-Pym	4-Pyr-CH ₂
1-2303	3,4-diF-Ph	4-Pym	3-Pyr-CH ₂
1-2304	3,4-diF-Ph	4-Pym	4-Pym-CH ₂
1-2305	3,4-diF-Ph	4-Pym	5-Pym-CH ₂
1-2306	3,4-diF-Ph	4-Pym	2-Pym-CH ₂
1-2307	3-Cl-Ph	4-Pym	H ₂ N-(CH ₂) ₃
1-2308	3-Cl-Ph	4-Pym	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-2309	3-Cl-Ph	4-Pym	EtNH-(CH ₂) ₃
1-2310	3-Cl-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
1-2311	3-Cl-Ph	4-Pym	1-Azt-(CH ₂) ₃
1-2312	3-Cl-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
1-2313	3-Cl-Ph	4-Pym	1-Pip-(CH ₂) ₃
1-2314	3-Cl-Ph	4-Pym	1-Mor-(CH ₂) ₃
1-2315	3-Cl-Ph	4-Pym	1-Tmor-(CH ₂) ₃
1-2316	3-Cl-Ph	4-Pym	1-Piz-(CH ₂) ₃
1-2317	3-Cl-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2318	3-Cl-Ph	4-Pym	3-Azt
1-2319	3-Cl-Ph	4-Pym	1-Me-3-Azt
1-2320	3-Cl-Ph	4-Pym	3-Pyrd
1-2321	3-Cl-Ph	4-Pym	1-Me-3-Pyrd
1-2322	3-Cl-Ph	4-Pym	4-Pip
1-2323	3-Cl-Ph	4-Pym	4-(3,4-deH-Pip)
1-2324	3-Cl-Ph	4-Pym	1-Me-4-Pip
1-2325	3-Cl-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
1-2326	3-Cl-Ph	4-Pym	1-Piz
1-2327	3-Cl-Ph	4-Pym	4-Me-1-Piz
1-2328	3-Cl-Ph	4-Pym	4-Pyr
1-2329	3-Cl-Ph	4-Pym	3-Pyr
1-2330	3-Cl-Ph	4-Pym	4-Pym
1-2331	3-Cl-Ph	4-Pym	5-Pym
1-2332	3-Cl-Ph	4-Pym	3-Pyrd-CH ₂
1-2333	3-Cl-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
1-2334	3-Cl-Ph	4-Pym	4-Pip-CH ₂
1-2335	3-Cl-Ph	4-Pym	1-Me-4-Pip-CH ₂
1-2336	3-Cl-Ph	4-Pym	2-Piz-CH ₂
1-2337	3-Cl-Ph	4-Pym	4-Pyr-CH ₂
1-2338	3-Cl-Ph	4-Pym	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
1-2339	3-Cl-Ph	4-Pym	4-Pym-CH ₂
1-2340	3-Cl-Ph	4-Pym	5-Pym-CH ₂
1-2341	3-Cl-Ph	4-Pym	2-Pym-CH ₂
1-2342	Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
1-2343	Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
1-2344	Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
1-2345	Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
1-2346	Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
1-2347	Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
1-2348	Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
1-2349	Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
1-2350	Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
1-2351	Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
1-2352	Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2353	Ph	2-MeO-4-Pym	3-Azt
1-2354	Ph	2-MeO-4-Pym	1-Me-3-Azt
1-2355	Ph	2-MeO-4-Pym	3-Pyrd
1-2356	Ph	2-MeO-4-Pym	1-Me-3-Pyrd
1-2357	Ph	2-MeO-4-Pym	4-Pip
1-2358	Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
1-2359	Ph	2-MeO-4-Pym	1-Me-4-Pip
1-2360	Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
1-2361	Ph	2-MeO-4-Pym	1-Piz
1-2362	Ph	2-MeO-4-Pym	4-Me-1-Piz
1-2363	Ph	2-MeO-4-Pym	4-Pyr
1-2364	Ph	2-MeO-4-Pym	3-Pyr
1-2365	Ph	2-MeO-4-Pym	4-Pym
1-2366	Ph	2-MeO-4-Pym	5-Pym
1-2367	Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
1-2368	Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
1-2369	Ph	2-MeO-4-Pym	4-Pip-CH ₂
1-2370	Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
1-2371	Ph	2-MeO-4-Pym	2-Piz-CH ₂
1-2372	Ph	2-MeO-4-Pym	4-Pyr-CH ₂
1-2373	Ph	2-MeO-4-Pym	3-Pyr-CH ₂
1-2374	Ph	2-MeO-4-Pym	4-Pym-CH ₂
1-2375	Ph	2-MeO-4-Pym	5-Pym-CH ₂
1-2376	Ph	2-MeO-4-Pym	2-Pym-CH ₂
1-2377	4-F-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
1-2378	4-F-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
1-2379	4-F-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
1-2380	4-F-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
1-2381	4-F-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
1-2382	4-F-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
1-2383	4-F-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
1-2384	4-F-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
1-2385	4-F-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
1-2386	4-F-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
1-2387	4-F-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2388	4-F-Ph	2-MeO-4-Pym	3-Azt
1-2389	4-F-Ph	2-MeO-4-Pym	1-Me-3-Azt
1-2390	4-F-Ph	2-MeO-4-Pym	3-Pyrd
1-2391	4-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
1-2392	4-F-Ph	2-MeO-4-Pym	4-Pip
1-2393	4-F-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
1-2394	4-F-Ph	2-MeO-4-Pym	1-Me-4-Pip
1-2395	4-F-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
1-2396	4-F-Ph	2-MeO-4-Pym	1-Piz
1-2397	4-F-Ph	2-MeO-4-Pym	4-Me-1-Piz
1-2398	4-F-Ph	2-MeO-4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
1-2399	4-F-Ph	2-MeO-4-Pym	3-Pyr
1-2400	4-F-Ph	2-MeO-4-Pym	4-Pym
1-2401	4-F-Ph	2-MeO-4-Pym	5-Pym
1-2402	4-F-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
1-2403	4-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
1-2404	4-F-Ph	2-MeO-4-Pym	4-Pip-CH ₂
1-2405	4-F-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
1-2406	4-F-Ph	2-MeO-4-Pym	2-Piz-CH ₂
1-2407	4-F-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
1-2408	4-F-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
1-2409	4-F-Ph	2-MeO-4-Pym	4-Pym-CH ₂
1-2410	4-F-Ph	2-MeO-4-Pym	5-Pym-CH ₂
1-2411	4-F-Ph	2-MeO-4-Pym	2-Pym-CH ₂
1-2412	3-F-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
1-2413	3-F-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
1-2414	3-F-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
1-2415	3-F-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₂
1-2416	3-F-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
1-2417	3-F-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
1-2418	3-F-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
1-2419	3-F-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
1-2420	3-F-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
1-2421	3-F-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
1-2422	3-F-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2423	3-F-Ph	2-MeO-4-Pym	3-Azt
1-2424	3-F-Ph	2-MeO-4-Pym	1-Me-3-Azt
1-2425	3-F-Ph	2-MeO-4-Pym	3-Pyrd
1-2426	3-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
1-2427	3-F-Ph	2-MeO-4-Pym	4-Pip
1-2428	3-F-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2429	3-F-Ph	2-MeO-4-Pym	1-Me-4-Pip
1-2430	3-F-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
1-2431	3-F-Ph	2-MeO-4-Pym	1-Piz
1-2432	3-F-Ph	2-MeO-4-Pym	4-Me-1-Piz
1-2433	3-F-Ph	2-MeO-4-Pym	4-Pyr
1-2434	3-F-Ph	2-MeO-4-Pym	3-Pyr
1-2435	3-F-Ph	2-MeO-4-Pym	4-Pym
1-2436	3-F-Ph	2-MeO-4-Pym	5-Pym
1-2437	3-F-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
1-2438	3-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
1-2439	3-F-Ph	2-MeO-4-Pym	4-Pip-CH ₂
1-2440	3-F-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
1-2441	3-F-Ph	2-MeO-4-Pym	2-Piz-CH ₂
1-2442	3-F-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
1-2443	3-F-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
1-2444	3-F-Ph	2-MeO-4-Pym	4-Pym-CH ₂
1-2445	3-F-Ph	2-MeO-4-Pym	5-Pym-CH ₂
1-2446	3-F-Ph	2-MeO-4-Pym	2-Pym-CH ₂
1-2447	3,4-diF-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₂
1-2448	3,4-diF-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
1-2449	3,4-diF-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
1-2450	3,4-diF-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
1-2451	3,4-diF-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
1-2452	3,4-diF-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
1-2453	3,4-diF-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
1-2454	3,4-diF-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
1-2455	3,4-diF-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
1-2456	3,4-diF-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
1-2457	3,4-diF-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2458	3,4-diF-Ph	2-MeO-4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
1-2459	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Azt
1-2460	3,4-diF-Ph	2-MeO-4-Pym	3-Pyrd
1-2461	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
1-2462	3,4-diF-Ph	2-MeO-4-Pym	4-Pip
1-2463	3,4-diF-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
1-2464	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-Pip
1-2465	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
1-2466	3,4-diF-Ph	2-MeO-4-Pym	1-Piz
1-2467	3,4-diF-Ph	2-MeO-4-Pym	4-Me-1-Piz
1-2468	3,4-diF-Ph	2-MeO-4-Pym	4-Pyr
1-2469	3,4-diF-Ph	2-MeO-4-Pym	3-Pyr
1-2470	3,4-diF-Ph	2-MeO-4-Pym	4-Pym
1-2471	3,4-diF-Ph	2-MeO-4-Pym	5-Pym
1-2472	3,4-diF-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
1-2473	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
1-2474	3,4-diF-Ph	2-MeO-4-Pym	4-Pip-CH ₂
1-2475	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
1-2476	3,4-diF-Ph	2-MeO-4-Pym	2-Piz-CH ₂
1-2477	3,4-diF-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
1-2478	3,4-diF-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
1-2479	3,4-diF-Ph	2-MeO-4-Pym	4-Pym-CH ₂
1-2480	3,4-diF-Ph	2-MeO-4-Pym	5-Pym-CH ₂
1-2481	3,4-diF-Ph	2-MeO-4-Pym	2-Pym-CH ₂
1-2482	3-Cl-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
1-2483	3-Cl-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
1-2484	3-Cl-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
1-2485	3-Cl-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
1-2486	3-Cl-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
1-2487	3-Cl-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
1-2488	3-Cl-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
1-2489	3-Cl-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
1-2490	3-Cl-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
1-2491	3-Cl-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
1-2492	3-Cl-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
1-2493	3-Cl-Ph	2-MeO-4-Pym	3-Azt
1-2494	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Azt
1-2495	3-Cl-Ph	2-MeO-4-Pym	3-Pyrd
1-2496	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
1-2497	3-Cl-Ph	2-MeO-4-Pym	4-Pip
1-2498	3-Cl-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
1-2499	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-Pip
1-2500	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
1-2501	3-Cl-Ph	2-MeO-4-Pym	1-Piz
1-2502	3-Cl-Ph	2-MeO-4-Pym	4-Me-1-Piz
1-2503	3-Cl-Ph	2-MeO-4-Pym	4-Pyr
1-2504	3-Cl-Ph	2-MeO-4-Pym	3-Pyr
1-2505	3-Cl-Ph	2-MeO-4-Pym	4-Pym
1-2506	3-Cl-Ph	2-MeO-4-Pym	5-Pym
1-2507	3-Cl-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
1-2508	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
1-2509	3-Cl-Ph	2-MeO-4-Pym	4-Pip-CH ₂
1-2510	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
1-2511	3-Cl-Ph	2-MeO-4-Pym	2-Piz-CH ₂
1-2512	3-Cl-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
1-2513	3-Cl-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
1-2514	3-Cl-Ph	2-MeO-4-Pym	4-Pym-CH ₂
1-2515	3-Cl-Ph	2-MeO-4-Pym	5-Pym-CH ₂
1-2516	3-Cl-Ph	2-MeO-4-Pym	2-Pym-CH ₂
1-2517	4-F-Ph	4-Pyr	H ₂ N-CH ₂ CH=CH
1-2518	4-F-Ph	4-Pyr	MeNH-CH ₂ CH=CH

Compound No.	R ¹	R ²	R ³
1-2519	4-F-Ph	4-Pyr	Me ₂ N-CH ₂ CH=CH
1-2520	4-F-Ph	4-Pyr	3-Pip-CH ₂
1-2521	4-F-Ph	4-Pyr	1-Me-3-Pip-CH ₂
1-2522	4-F-Ph	4-Pyr	2-Me-4-Pip
1-2523	4-F-Ph	4-Pyr	2,2,6,6-tetraMe-4-Pip
1-2524	4-F-Ph	4-Pyr	1-Ac-4-Pip
1-2525	4-F-Ph	4-Pyr	1-Ac-4-(3,4-deH-Pip)
1-2526	4-F-Ph	4-Pyr	4-OH-4-Pip
1-2527	4-F-Ph	4-Pyr	4-OH-1-Me-4-Pip
1-2528	4-F-Ph	4-Pyr	AcNH-(CH ₂) ₃
1-2529	4-F-Ph	4-Pyr	4-NH ₂ -cHx
1-2530	4-F-Ph	4-Pyr	4-Pyr-CH(OH)
1-2531	4-F-Ph	4-Pyr	3-Pyr-CH(OH)
1-2532	4-F-Ph	4-Pyr	2-Pyr-CH(OH)
1-2533	4-F-Ph	4-Pyr	CF ₃ CONH-(CH ₂) ₃
1-2534	4-F-Ph	4-Pyr	BzNH-(CH ₂) ₃
1-2535	4-F-Ph	4-Pyr	2,4,6-triF-BzNH-CH ₂
1-2536	4-F-Ph	4-Pyr	MeSO ₂ NH-(CH ₂) ₃
1-2537	4-F-Ph	4-Pyr	1-NO ₂ (CH ₂) ₂ -4-Pip
1-2538	4-F-Ph	4-Pyr	2,3,5,6-tetraF-4-Pyr
1-2539	4-F-Ph	4-Pyr	3-Qun
1-2540	4-F-Ph	4-Pyr	3-(2,3-deH-Qun)
1-2541	4-F-Ph	4-Pyr	3-ABO
1-2542	4-F-Ph	4-Pyr	8-Me-3-ABO
1-2543	4-F-Ph	4-Pyr	3-(2,3-deH-ABO)
1-2544	4-F-Ph	4-Pyr	8-Me-3-(2,3-deH-ABO)
1-2545	4-F-Ph	4-Pyr	3-ABN
1-2546	4-F-Ph	4-Pyr	9-Me-3-ABN
1-2547	4-F-Ph	4-Pyr	3-(2,3-deH-ABN)
1-2548	4-F-Ph	4-Pyr	9-Me-3-(2,3-deH-ABN)

Compound No.	R ¹	R ²	R ³
1-2549	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂ CH=CH
1-2550	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂ CH=CH
1-2551	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂ CH=CH
1-2552	4-F-Ph	2-NH ₂ -4-Pym	3-Pip-CH ₂
1-2553	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pip-CH ₂
1-2554	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-Pip
1-2555	4-F-Ph	2-NH ₂ -4-Pym	2,2,6,6-tetraMe-4-Pip
1-2556	4-F-Ph	2-NH ₂ -4-Pym	1-Ac-4-Pip
1-2557	4-F-Ph	2-NH ₂ -4-Pym	1-Ac-4-(3,4-deH-Pip)
1-2558	4-F-Ph	2-NH ₂ -4-Pym	4-OH-4-Pip
1-2559	4-F-Ph	2-NH ₂ -4-Pym	4-OH-1-Me-4-Pip
1-2560	4-F-Ph	2-NH ₂ -4-Pym	AcNH-(CH ₂) ₃
1-2561	4-F-Ph	2-NH ₂ -4-Pym	4-NH ₂ -cHx
1-2562	4-F-Ph	2-NH ₂ -4-Pym	3-Qun
1-2563	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-Qun)
1-2564	4-F-Ph	2-NH ₂ -4-Pym	3-ABO
1-2565	4-F-Ph	2-NH ₂ -4-Pym	8-Me-3-ABO
1-2566	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-ABO)
1-2567	4-F-Ph	2-NH ₂ -4-Pym	8-Me-3-(2,3-deH-ABO)
1-2568	4-F-Ph	2-NH ₂ -4-Pym	3-ABN
1-2569	4-F-Ph	2-NH ₂ -4-Pym	9-Me-3-ABN
1-2570	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-ABN)
1-2571	4-F-Ph	2-NH ₂ -4-Pym	9-Me-3-(2,3-deH-ABN)
1-2572	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂ CH=CH
1-2573	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂ CH=CH
1-2574	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂ CH=CH
1-2575	4-F-Ph	2-MeNH-4-Pym	3-Pip-CH ₂
1-2576	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pip-CH ₂
1-2577	4-F-Ph	2-MeNH-4-Pym	2-Me-4-Pip
1-2578	4-F-Ph	2-MeNH-4-Pym	2,2,6,6-tetraMe-4-Pip

Compound No.	R ¹	R ²	R ³
1-2579	4-F-Ph	2-MeNH-4-Pym	1-Ac-4-Pip
1-2580	4-F-Ph	2-MeNH-4-Pym	1-Ac-4-(3,4-deH-Pip)
1-2581	4-F-Ph	2-MeNH-4-Pym	4-OH-4-Pip
1-2582	4-F-Ph	2-MeNH-4-Pym	4-OH-1-Me-4-Pip
1-2583	4-F-Ph	2-MeNH-4-Pym	AcNH-(CH ₂) ₃
1-2584	4-F-Ph	2-MeNH-4-Pym	4-NH ₂ -cHx
1-2585	4-F-Ph	2-MeNH-4-Pym	3-Qun
1-2586	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-Qun)
1-2587	4-F-Ph	2-MeNH-4-Pym	3-ABO
1-2588	4-F-Ph	2-MeNH-4-Pym	8-Me-3-ABO
1-2589	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-ABO)
1-2590	4-F-Ph	2-MeNH-4-Pym	8-Me-3-(2,3-deH-ABO)
1-2591	4-F-Ph	2-MeNH-4-Pym	3-ABN
1-2592	4-F-Ph	2-MeNH-4-Pym	9-Me-3-ABN
1-2593	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-ABN)
1-2594	4-F-Ph	2-MeNH-4-Pym	9-Me-3-(2,3-deH-ABN)
1-2595	4-F-Ph	2-BnNH-4-Pyr	4-Pip
1-2596	4-F-Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
1-2597	4-F-Ph	2-BnNH-4-Pym	4-Pip
1-2598	4-F-Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)
1-2599	4-F-Ph	2-(α -Me-BnNH)-4-Pyr	4-Pip
1-2600	4-F-Ph	2-(α -Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)
1-2601	4-F-Ph	2-(β -Me-BnNH)-4-Pym	4-Pip
1-2602	4-F-Ph	2-(β -Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
1-2603	3-Cl-Ph	2-BnNH-4-Pyr	4-Pip
1-2604	3-Cl-Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
1-2605	3-Cl-Ph	2-BnNH-4-Pym	4-Pip
1-2606	3-Cl-Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)
1-2607	3-Cl-Ph	2-(β -Me-BnNH)-4-Pyr	4-Pip
1-2608	3-Cl-Ph	2-(β -Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2609	3-Cl-Ph	2-(-Me-BnNH)-4-Pym	4-Pip
1-2610	3-Cl-Ph	2-(-Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
1-2611	3-CF ₃ -Ph	2-BnNH-4-Pyr	4-Pip
1-2612	3-CF ₃ -Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
1-2613	3-CF ₃ -Ph	2-BnNH-4-Pym	4-Pip
1-2614	3-CF ₃ -Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)
1-2615	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pyr	4-Pip
1-2616	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)
1-2617	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pym	4-Pip
1-2618	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
1-2619	4-F-Ph	4-Pyr	2-NH ₂ -4-Pym
1-2620	4-F-Ph	4-Pyr	2-MeNH-4-Pym
1-2621	4-F-Ph	4-Pyr	2-NH ₂ -4-Pyr
1-2622	4-F-Ph	4-Pyr	2-MeNH-4-Pyr
1-2623	4-F-Ph	4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2624	4-F-Ph	4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
1-2625	4-F-Ph	4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
1-2626	4-F-Ph	4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2627	4-F-Ph	4-Pyr	3-(3,4-deH-Pip)
1-2628	4-F-Ph	4-Pyr	1-Me-3-(3,4-deH-Pip)
1-2629	4-F-Ph	4-Pyr	1-Et-4-(3,4-deH-Pip)
1-2630	4-F-Ph	4-Pyr	1-Pr-4-(3,4-deH-Pip)
1-2631	4-F-Ph	4-Pyr	1-Pr-4-Pip
1-2632	4-F-Ph	4-Pyr	1-iPr-4-(3,4-deH-Pip)
1-2633	4-F-Ph	4-Pyr	1-iPr-4-Pip
1-2634	4-F-Ph	4-Pyr	1-Bu-4-(3,4-deH-Pip)
1-2635	4-F-Ph	4-Pyr	1-tBu-4-(3,4-deH-Pip)
1-2636	4-F-Ph	4-Pyr	1-Pn-4-(3,4-deH-Pip)
1-2637	4-F-Ph	4-Pyr	1-Hx-4-(3,4-deH-Pip)
1-2638	4-F-Ph	4-Pyr	1-Hp-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2639	4-F-Ph	4-Pyr	1-Oc-4-(3,4-deH-Pip)
1-2640	4-F-Ph	4-Pyr	1-Nn-4-(3,4-deH-Pip)
1-2641	4-F-Ph	4-Pyr	1-cPr-4-(3,4-deH-Pip)
1-2642	4-F-Ph	4-Pyr	1-cPn-4-(3,4-deH-Pip)
1-2643	4-F-Ph	4-Pyr	1-cHx-4-(3,4-deH-Pip)
1-2644	4-F-Ph	4-Pyr	1-Bn-4-(3,4-deH-Pip)
1-2645	4-F-Ph	4-Pyr	1-Phet-4-(3,4-deH-Pip)
1-2646	4-F-Ph	4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
1-2647	4-F-Ph	4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
1-2648	4-F-Ph	4-Pyr	1-Allyl-4-(3,4-deH-Pip)
1-2649	4-F-Ph	4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
1-2650	4-F-Ph	4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
1-2651	4-F-Ph	4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
1-2652	4-F-Ph	4-Pyr	1,2,2,6,6-pentaMe-4-Pip
1-2653	4-F-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
1-2654	4-F-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
1-2655	4-F-Ph	4-Pyr	7-octaH-Ind
1-2656	4-F-Ph	4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
1-2657	4-F-Ph	4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
1-2658	4-F-Ph	4-Pyr	8-octaH-Qui
1-2659	4-F-Ph	4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
1-2660	4-F-Ph	4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
1-2661	4-F-Ph	4-Pyr	2,2-diMe-4-(4,5-deH-Pip)
1-2662	4-F-Ph	4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)
1-2663	4-F-Ph	4-Pyr	2,6-diMe-4-(3,4-deH-Pip)
1-2664	4-F-Ph	4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
1-2665	4-F-Ph	4-Pyr	2-Me-4-(3,4-deH-Pip)
1-2666	4-F-Ph	4-Pyr	1,2-diMe-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2667	4-F-Ph	4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
1-2668	4-F-Ph	4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
1-2669	4-F-Ph	4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
1-2670	4-F-Ph	4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)
1-2671	4-F-Ph	4-Pyr	2-Et-4-(3,4-deH-Pip)
1-2672	4-F-Ph	4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
1-2673	4-F-Ph	4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
1-2674	4-F-Ph	4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
1-2675	4-F-Ph	4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
1-2676	4-F-Ph	4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
1-2677	4-F-Ph	4-Pyr	2-Pr-4-(3,4-deH-Pip)
1-2678	4-F-Ph	4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
1-2679	4-F-Ph	4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
1-2680	4-F-Ph	4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
1-2681	4-F-Ph	4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
1-2682	4-F-Ph	4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
1-2683	4-F-Ph	4-Pyr	2-Bu-4-(3,4-deH-Pip)
1-2684	4-F-Ph	4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
1-2685	4-F-Ph	4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
1-2686	4-F-Ph	4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
1-2687	4-F-Ph	4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
1-2688	4-F-Ph	4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
1-2689	4-F-Ph	4-Pyr	2-Allyl-4-(3,4-deH-Pip)
1-2690	4-F-Ph	4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
1-2691	4-F-Ph	4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)
1-2692	4-F-Ph	4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)
1-2693	4-F-Ph	4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
1-2694	4-F-Ph	4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
1-2695	4-F-Ph	4-Pyr	2-Bn-4-(3,4-deH-Pip)
1-2696	4-F-Ph	4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2697	4-F-Ph	4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
1-2698	4-F-Ph	4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
1-2699	4-F-Ph	4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
1-2700	4-F-Ph	4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)
1-2701	4-F-Ph	4-Pyr	2-Phet-4-(3,4-deH-Pip)
1-2702	4-F-Ph	4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
1-2703	4-F-Ph	4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
1-2704	4-F-Ph	4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
1-2705	4-F-Ph	4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
1-2706	4-F-Ph	4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
1-2707	4-F-Ph	4-Pyr	2-Me-4-(4,5-deH-Pip)
1-2708	4-F-Ph	4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
1-2709	4-F-Ph	4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
1-2710	4-F-Ph	4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
1-2711	4-F-Ph	4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
1-2712	4-F-Ph	4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
1-2713	4-F-Ph	4-Pyr	2-Et-4-(4,5-deH-Pip)
1-2714	4-F-Ph	4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
1-2715	4-F-Ph	4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
1-2716	4-F-Ph	4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
1-2717	4-F-Ph	4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
1-2718	4-F-Ph	4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
1-2719	4-F-Ph	4-Pyr	2-Pr-4-(4,5-deH-Pip)
1-2720	4-F-Ph	4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
1-2721	4-F-Ph	4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)
1-2722	4-F-Ph	4-Pyr	1,2-diPr-4-(4,5-deH-Pip)
1-2723	4-F-Ph	4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
1-2724	4-F-Ph	4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
1-2725	4-F-Ph	4-Pyr	2-Bu-4-(4,5-deH-Pip)
1-2726	4-F-Ph	4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2727	4-F-Ph	4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
1-2728	4-F-Ph	4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
1-2729	4-F-Ph	4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
1-2730	4-F-Ph	4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)
1-2731	4-F-Ph	4-Pyr	2-Allyl-4-(4,5-deH-Pip)
1-2732	4-F-Ph	4-Pyr	1-Me-2-Allyl-4-(4,5-deH-Pip)
1-2733	4-F-Ph	4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
1-2734	4-F-Ph	4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
1-2735	4-F-Ph	4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
1-2736	4-F-Ph	4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
1-2737	4-F-Ph	4-Pyr	2-Bn-4-(4,5-deH-Pip)
1-2738	4-F-Ph	4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
1-2739	4-F-Ph	4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
1-2740	4-F-Ph	4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
1-2741	4-F-Ph	4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
1-2742	4-F-Ph	4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
1-2743	4-F-Ph	4-Pyr	2-Phet-4-(4,5-deH-Pip)
1-2744	4-F-Ph	4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
1-2745	4-F-Ph	4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
1-2746	4-F-Ph	4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
1-2747	4-F-Ph	4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
1-2748	4-F-Ph	4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
1-2749	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2750	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂ C(Me) ₂ CH ₂
1-2751	4-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂ C(Me) ₂ CH ₂
1-2752	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2753	4-F-Ph	2-NH ₂ -4-Pym	3-(3,4-deH-Pip)
1-2754	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-(3,4-deH-Pip)
1-2755	4-F-Ph	2-NH ₂ -4-Pym	1-Et-4-(3,4-deH-Pip)
1-2756	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2757	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-4-Pip
1-2758	4-F-Ph	2-NH ₂ -4-Pym	1-iPr-4-(3,4-deH-Pip)
1-2759	4-F-Ph	2-NH ₂ -4-Pym	1-iPr-4-Pip
1-2760	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-4-(3,4-deH-Pip)
1-2761	4-F-Ph	2-NH ₂ -4-Pym	1-tBu-4-(3,4-deH-Pip)
1-2762	4-F-Ph	2-NH ₂ -4-Pym	1-Pn-4-(3,4-deH-Pip)
1-2763	4-F-Ph	2-NH ₂ -4-Pym	1-Hx-4-(3,4-deH-Pip)
1-2764	4-F-Ph	2-NH ₂ -4-Pym	1-Hp-4-(3,4-deH-Pip)
1-2765	4-F-Ph	2-NH ₂ -4-Pym	1-Oc-4-(3,4-deH-Pip)
1-2766	4-F-Ph	2-NH ₂ -4-Pym	1-Nn-4-(3,4-deH-Pip)
1-2767	4-F-Ph	2-NH ₂ -4-Pym	1-cPr-4-(3,4-deH-Pip)
1-2768	4-F-Ph	2-NH ₂ -4-Pym	1-cPn-4-(3,4-deH-Pip)
1-2769	4-F-Ph	2-NH ₂ -4-Pym	1-cHx-4-(3,4-deH-Pip)
1-2770	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-(3,4-deH-Pip)
1-2771	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-4-(3,4-deH-Pip)
1-2772	4-F-Ph	2-NH ₂ -4-Pym	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
1-2773	4-F-Ph	2-NH ₂ -4-Pym	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
1-2774	4-F-Ph	2-NH ₂ -4-Pym	1-Allyl-4-(3,4-deH-Pip)
1-2775	4-F-Ph	2-NH ₂ -4-Pym	1-Propargyl-4-(3,4-deH-Pip)
1-2776	4-F-Ph	2-NH ₂ -4-Pym	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
1-2777	4-F-Ph	2-NH ₂ -4-Pym	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
1-2778	4-F-Ph	2-NH ₂ -4-Pym	1,2,2,6,6-pentaMe-4-Pip
1-2779	4-F-Ph	2-NH ₂ -4-Pym	7-(1,2,3,5,6,8a-hexaH-Ind)
1-2780	4-F-Ph	2-NH ₂ -4-Pym	7-(1,2,3,5,8,8a-hexaH-Ind)
1-2781	4-F-Ph	2-NH ₂ -4-Pym	7-octaH-Ind
1-2782	4-F-Ph	2-NH ₂ -4-Pym	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
1-2783	4-F-Ph	2-NH ₂ -4-Pym	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
1-2784	4-F-Ph	2-NH ₂ -4-Pym	8-octaH-Qui

Compound No.	R ¹	R ²	R ³
1-2785	4-F-Ph	2-NH ₂ -4-Pym	2,2-diMe-4-(3,4-deH-Pip)
1-2786	4-F-Ph	2-NH ₂ -4-Pym	1,2,2-triMe-4-(3,4-deH-Pip)
1-2787	4-F-Ph	2-NH ₂ -4-Pym	2,2-diMe-4-(4,5-deH-Pip)
1-2788	4-F-Ph	2-NH ₂ -4-Pym	1,2,2-triMe-4-(4,5-deH-Pip)
1-2789	4-F-Ph	2-NH ₂ -4-Pym	2,6-diMe-4-(3,4-deH-Pip)
1-2790	4-F-Ph	2-NH ₂ -4-Pym	1,2,6-triMe-4-(3,4-deH-Pip)
1-2791	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-(3,4-deH-Pip)
1-2792	4-F-Ph	2-NH ₂ -4-Pym	1,2-diMe-4-(3,4-deH-Pip)
1-2793	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Me-4-(3,4-deH-Pip)
1-2794	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Me-4-(3,4-deH-Pip)
1-2795	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Me-4-(3,4-deH-Pip)
1-2796	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Me-4-(3,4-deH-Pip)
1-2797	4-F-Ph	2-NH ₂ -4-Pym	2-Et-4-(3,4-deH-Pip)
1-2798	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Et-4-(3,4-deH-Pip)
1-2799	4-F-Ph	2-NH ₂ -4-Pym	1,2-diEt-4-(3,4-deH-Pip)
1-2800	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Et-4-(3,4-deH-Pip)
1-2801	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Et-4-(3,4-deH-Pip)
1-2802	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Et-4-(3,4-deH-Pip)
1-2803	4-F-Ph	2-NH ₂ -4-Pym	2-Pr-4-(3,4-deH-Pip)
1-2804	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Pr-4-(3,4-deH-Pip)
1-2805	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Pr-4-(3,4-deH-Pip)
1-2806	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPr-4-(3,4-deH-Pip)
1-2807	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Pr-4-(3,4-deH-Pip)
1-2808	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Pr-4-(3,4-deH-Pip)
1-2809	4-F-Ph	2-NH ₂ -4-Pym	2-Bu-4-(3,4-deH-Pip)
1-2810	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bu-4-(3,4-deH-Pip)
1-2811	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bu-4-(3,4-deH-Pip)
1-2812	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bu-4-(3,4-deH-Pip)
1-2813	4-F-Ph	2-NH ₂ -4-Pym	1,2-diBu-4-(3,4-deH-Pip)
1-2814	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bu-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2815	4-F-Ph	2-NH ₂ -4-Pym	2-Allyl-4-(3,4-deH-Pip)
1-2816	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Allyl-4-(3,4-deH-Pip)
1-2817	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Allyl-4-(3,4-deH-Pip)
1-2818	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Allyl-4-(3,4-deH-Pip)
1-2819	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Allyl-4-(3,4-deH-Pip)
1-2820	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Allyl-4-(3,4-deH-Pip)
1-2821	4-F-Ph	2-NH ₂ -4-Pym	2-Bn-4-(3,4-deH-Pip)
1-2822	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bn-4-(3,4-deH-Pip)
1-2823	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bn-4-(3,4-deH-Pip)
1-2824	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bn-4-(3,4-deH-Pip)
1-2825	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Bn-4-(3,4-deH-Pip)
1-2826	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bn-4-(3,4-deH-Pip)
1-2827	4-F-Ph	2-NH ₂ -4-Pym	2-Phet-4-(3,4-deH-Pip)
1-2828	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Phet-4-(3,4-deH-Pip)
1-2829	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Phet-4-(3,4-deH-Pip)
1-2830	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Phet-4-(3,4-deH-Pip)
1-2831	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Phet-4-(3,4-deH-Pip)
1-2832	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPhet-4-(3,4-deH-Pip)
1-2833	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-(4,5-deH-Pip)
1-2834	4-F-Ph	2-NH ₂ -4-Pym	1,2-diMe-4-(4,5-deH-Pip)
1-2835	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Me-4-(4,5-deH-Pip)
1-2836	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Me-4-(4,5-deH-Pip)
1-2837	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Me-4-(4,5-deH-Pip)
1-2838	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Me-4-(4,5-deH-Pip)
1-2839	4-F-Ph	2-NH ₂ -4-Pym	2-Et-4-(4,5-deH-Pip)
1-2840	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Et-4-(4,5-deH-Pip)
1-2841	4-F-Ph	2-NH ₂ -4-Pym	1,2-diEt-4-(4,5-deH-Pip)
1-2842	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Et-4-(4,5-deH-Pip)
1-2843	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Et-4-(4,5-deH-Pip)
1-2844	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Et-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2845	4-F-Ph	2-NH ₂ -4-Pym	2-Pr-4-(4,5-deH-Pip)
1-2846	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Pr-4-(4,5-deH-Pip)
1-2847	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Pr-4-(4,5-deH-Pip)
1-2848	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPr-4-(4,5-deH-Pip)
1-2849	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Pr-4-(4,5-deH-Pip)
1-2850	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Pr-4-(4,5-deH-Pip)
1-2851	4-F-Ph	2-NH ₂ -4-Pym	2-Bu-4-(4,5-deH-Pip)
1-2852	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bu-4-(4,5-deH-Pip)
1-2853	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bu-4-(4,5-deH-Pip)
1-2854	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bu-4-(4,5-deH-Pip)
1-2855	4-F-Ph	2-NH ₂ -4-Pym	1,2-diBu-4-(4,5-deH-Pip)
1-2856	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bu-4-(4,5-deH-Pip)
1-2857	4-F-Ph	2-NH ₂ -4-Pym	2-Allyl-4-(4,5-deH-Pip)
1-2858	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Allyl-4-(4,5-deH-Pip)
1-2859	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Allyl-4-(4,5-deH-Pip)
1-2860	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Allyl-4-(4,5-deH-Pip)
1-2861	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Allyl-4-(4,5-deH-Pip)
1-2862	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Allyl-4-(4,5-deH-Pip)
1-2863	4-F-Ph	2-NH ₂ -4-Pym	2-Bn-4-(4,5-deH-Pip)
1-2864	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bn-4-(4,5-deH-Pip)
1-2865	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bn-4-(4,5-deH-Pip)
1-2866	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bn-4-(4,5-deH-Pip)
1-2867	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Bn-4-(4,5-deH-Pip)
1-2868	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bn-4-(4,5-deH-Pip)
1-2869	4-F-Ph	2-NH ₂ -4-Pym	2-Phet-4-(4,5-deH-Pip)
1-2870	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Phet-4-(4,5-deH-Pip)
1-2871	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Phet-4-(4,5-deH-Pip)
1-2872	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Phet-4-(4,5-deH-Pip)
1-2873	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Phet-4-(4,5-deH-Pip)
1-2874	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPhet-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2875	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2876	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂ C(Me) ₂ CH ₂
1-2877	4-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂ C(Me) ₂ CH ₂
1-2878	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
1-2879	4-F-Ph	2-MeNH-4-Pym	3-(3,4-deH-Pip)
1-2880	4-F-Ph	2-MeNH-4-Pym	1-Me-3-(3,4-deH-Pip)
1-2881	4-F-Ph	2-MeNH-4-Pym	1-Et-4-(3,4-deH-Pip)
1-2882	4-F-Ph	2-MeNH-4-Pym	1-Pr-4-(3,4-deH-Pip)
1-2883	4-F-Ph	2-MeNH-4-Pym	1-Pr-4-Pip
1-2884	4-F-Ph	2-MeNH-4-Pym	1-iPr-4-(3,4-deH-Pip)
1-2885	4-F-Ph	2-MeNH-4-Pym	1-iPr-4-Pip
1-2886	4-F-Ph	2-MeNH-4-Pym	1-Bu-4-(3,4-deH-Pip)
1-2887	4-F-Ph	2-MeNH-4-Pym	1-tBu-4-(3,4-deH-Pip)
1-2888	4-F-Ph	2-MeNH-4-Pym	1-Pn-4-(3,4-deH-Pip)
1-2889	4-F-Ph	2-MeNH-4-Pym	1-Hx-4-(3,4-deH-Pip)
1-2890	4-F-Ph	2-MeNH-4-Pym	1-Hp-4-(3,4-deH-Pip)
1-2891	4-F-Ph	2-MeNH-4-Pym	1-Oc-4-(3,4-deH-Pip)
1-2892	4-F-Ph	2-MeNH-4-Pym	1-Nn-4-(3,4-deH-Pip)
1-2893	4-F-Ph	2-MeNH-4-Pym	1-cPr-4-(3,4-deH-Pip)
1-2894	4-F-Ph	2-MeNH-4-Pym	1-cPn-4-(3,4-deH-Pip)
1-2895	4-F-Ph	2-MeNH-4-Pym	1-cHx-4-(3,4-deH-Pip)
1-2896	4-F-Ph	2-MeNH-4-Pym	1-Bn-4-(3,4-deH-Pip)
1-2897	4-F-Ph	2-MeNH-4-Pym	1-Phet-4-(3,4-deH-Pip)
1-2898	4-F-Ph	2-MeNH-4-Pym	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
1-2899	4-F-Ph	2-MeNH-4-Pym	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
1-2900	4-F-Ph	2-MeNH-4-Pym	1-Allyl-4-(3,4-deH-Pip)
1-2901	4-F-Ph	2-MeNH-4-Pym	1-Propargyl-4-(3,4-deH-Pip)
1-2902	4-F-Ph	2-MeNH-4-Pym	2,2,6,6-tetraMe-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2903	4-F-Ph	2-MeNH-4-Pym	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
1-2904	4-F-Ph	2-MeNH-4-Pym	1,2,2,6,6-pentaMe-4-Pip
1-2905	4-F-Ph	2-MeNH-4-Pym	7-(1,2,3,5,6,8a-hexaH-Ind)
1-2906	4-F-Ph	2-MeNH-4-Pym	7-(1,2,3,5,8,8a-hexaH-Ind)
1-2907	4-F-Ph	2-MeNH-4-Pym	7-octaH-Ind
1-2908	4-F-Ph	2-MeNH-4-Pym	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
1-2909	4-F-Ph	2-MeNH-4-Pym	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
1-2910	4-F-Ph	2-MeNH-4-Pym	8-octaH-Qui
1-2911	4-F-Ph	2-MeNH-4-Pym	2,2-diMe-4-(3,4-deH-Pip)
1-2912	4-F-Ph	2-MeNH-4-Pym	1,2,2-triMe-4-(3,4-deH-Pip)
1-2913	4-F-Ph	2-MeNH-4-Pym	2,2-diMe-4-(4,5-deH-Pip)
1-2914	4-F-Ph	2-MeNH-4-Pym	1,2,2-triMe-4-(4,5-deH-Pip)
1-2915	4-F-Ph	2-MeNH-4-Pym	2,6-diMe-4-(3,4-deH-Pip)
1-2916	4-F-Ph	2-MeNH-4-Pym	1,2,6-triMe-4-(3,4-deH-Pip)
1-2917	4-F-Ph	2-MeNH-4-Pym	2-Me-4-(3,4-deH-Pip)
1-2918	4-F-Ph	2-MeNH-4-Pym	1,2-diMe-4-(3,4-deH-Pip)
1-2919	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Me-4-(3,4-deH-Pip)
1-2920	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Me-4-(3,4-deH-Pip)
1-2921	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Me-4-(3,4-deH-Pip)
1-2922	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Me-4-(3,4-deH-Pip)
1-2923	4-F-Ph	2-MeNH-4-Pym	2-Et-4-(3,4-deH-Pip)
1-2924	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Et-4-(3,4-deH-Pip)
1-2925	4-F-Ph	2-MeNH-4-Pym	1,2-diEt-4-(3,4-deH-Pip)
1-2926	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Et-4-(3,4-deH-Pip)
1-2927	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Et-4-(3,4-deH-Pip)
1-2928	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Et-4-(3,4-deH-Pip)
1-2929	4-F-Ph	2-MeNH-4-Pym	2-Pr-4-(3,4-deH-Pip)
1-2930	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Pr-4-(3,4-deH-Pip)
1-2931	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Pr-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2932	4-F-Ph	2-MeNH-4-Pym	1,2-diPr-4-(3,4-deH-Pip)
1-2933	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Pr-4-(3,4-deH-Pip)
1-2934	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Pr-4-(3,4-deH-Pip)
1-2935	4-F-Ph	2-MeNH-4-Pym	2-Bu-4-(3,4-deH-Pip)
1-2936	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bu-4-(3,4-deH-Pip)
1-2937	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bu-4-(3,4-deH-Pip)
1-2938	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bu-4-(3,4-deH-Pip)
1-2939	4-F-Ph	2-MeNH-4-Pym	1,2-diBu-4-(3,4-deH-Pip)
1-2940	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bu-4-(3,4-deH-Pip)
1-2941	4-F-Ph	2-MeNH-4-Pym	2-Allyl-4-(3,4-deH-Pip)
1-2942	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Allyl-4-(3,4-deH-Pip)
1-2943	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Allyl-4-(3,4-deH-Pip)
1-2944	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Allyl-4-(3,4-deH-Pip)
1-2945	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Allyl-4-(3,4-deH-Pip)
1-2946	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Allyl-4-(3,4-deH-Pip)
1-2947	4-F-Ph	2-MeNH-4-Pym	2-Bn-4-(3,4-deH-Pip)
1-2948	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bn-4-(3,4-deH-Pip)
1-2949	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bn-4-(3,4-deH-Pip)
1-2950	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bn-4-(3,4-deH-Pip)
1-2951	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Bn-4-(3,4-deH-Pip)
1-2952	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bn-4-(3,4-deH-Pip)
1-2953	4-F-Ph	2-MeNH-4-Pym	2-Phet-4-(3,4-deH-Pip)
1-2954	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Phet-4-(3,4-deH-Pip)
1-2955	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Phet-4-(3,4-deH-Pip)
1-2956	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Phet-4-(3,4-deH-Pip)
1-2957	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Phet-4-(3,4-deH-Pip)
1-2958	4-F-Ph	2-MeNH-4-Pym	1,2-diPhet-4-(3,4-deH-Pip)
1-2959	4-F-Ph	2-MeNH-4-Pym	2-Me-4-(4,5-deH-Pip)
1-2960	4-F-Ph	2-MeNH-4-Pym	1,2-diMe-4-(4,5-deH-Pip)
1-2961	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Me-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2962	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Me-4-(4,5-deH-Pip)
1-2963	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Me-4-(4,5-deH-Pip)
1-2964	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Me-4-(4,5-deH-Pip)
1-2965	4-F-Ph	2-MeNH-4-Pym	2-Et-4-(4,5-deH-Pip)
1-2966	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Et-4-(4,5-deH-Pip)
1-2967	4-F-Ph	2-MeNH-4-Pym	1,2-diEt-4-(4,5-deH-Pip)
1-2968	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Et-4-(4,5-deH-Pip)
1-2969	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Et-4-(4,5-deH-Pip)
1-2970	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Et-4-(4,5-deH-Pip)
1-2971	4-F-Ph	2-MeNH-4-Pym	2-Pr-4-(4,5-deH-Pip)
1-2972	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Pr-4-(4,5-deH-Pip)
1-2973	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Pr-4-(4,5-deH-Pip)
1-2974	4-F-Ph	2-MeNH-4-Pym	1,2-diPr-4-(4,5-deH-Pip)
1-2975	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Pr-4-(4,5-deH-Pip)
1-2976	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Pr-4-(4,5-deH-Pip)
1-2977	4-F-Ph	2-MeNH-4-Pym	2-Bu-4-(4,5-deH-Pip)
1-2978	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bu-4-(4,5-deH-Pip)
1-2979	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bu-4-(4,5-deH-Pip)
1-2980	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bu-4-(4,5-deH-Pip)
1-2981	4-F-Ph	2-MeNH-4-Pym	1,2-diBu-4-(4,5-deH-Pip)
1-2982	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bu-4-(4,5-deH-Pip)
1-2983	4-F-Ph	2-MeNH-4-Pym	2-Allyl-4-(4,5-deH-Pip)
1-2984	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Allyl-4-(4,5-deH-Pip)
1-2985	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Allyl-4-(4,5-deH-Pip)
1-2986	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Allyl-4-(4,5-deH-Pip)
1-2987	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Allyl-4-(4,5-deH-Pip)
1-2988	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Allyl-4-(4,5-deH-Pip)
1-2989	4-F-Ph	2-MeNH-4-Pym	2-Bn-4-(4,5-deH-Pip)
1-2990	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bn-4-(4,5-deH-Pip)
1-2991	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bn-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-2992	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bn-4-(4,5-deH-Pip)
1-2993	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Bn-4-(4,5-deH-Pip)
1-2994	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bn-4-(4,5-deH-Pip)
1-2995	4-F-Ph	2-MeNH-4-Pym	2-Phet-4-(4,5-deH-Pip)
1-2996	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Phet-4-(4,5-deH-Pip)
1-2997	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Phet-4-(4,5-deH-Pip)
1-2998	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Phet-4-(4,5-deH-Pip)
1-2999	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Phet-4-(4,5-deH-Pip)
1-3000	4-F-Ph	2-MeNH-4-Pym	1,2-diPhet-4-(4,5-deH-Pip)
1-3001	4-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
1-3002	4-F-Ph	2-NH ₂ -4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
1-3003	4-F-Ph	2-NH ₂ -4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
1-3004	4-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
1-3005	4-F-Ph	2-NH ₂ -4-Pyr	3-(3,4-deH-Pip)
1-3006	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-(3,4-deH-Pip)
1-3007	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-4-(3,4-deH-Pip)
1-3008	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-4-(3,4-deH-Pip)
1-3009	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-4-Pip
1-3010	4-F-Ph	2-NH ₂ -4-Pyr	1-iPr-4-(3,4-deH-Pip)
1-3011	4-F-Ph	2-NH ₂ -4-Pyr	1-iPr-4-Pip
1-3012	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-4-(3,4-deH-Pip)
1-3013	4-F-Ph	2-NH ₂ -4-Pyr	1-tBu-4-(3,4-deH-Pip)
1-3014	4-F-Ph	2-NH ₂ -4-Pyr	1-Pn-4-(3,4-deH-Pip)
1-3015	4-F-Ph	2-NH ₂ -4-Pyr	1-Hx-4-(3,4-deH-Pip)
1-3016	4-F-Ph	2-NH ₂ -4-Pyr	1-Hp-4-(3,4-deH-Pip)
1-3017	4-F-Ph	2-NH ₂ -4-Pyr	1-Oc-4-(3,4-deH-Pip)
1-3018	4-F-Ph	2-NH ₂ -4-Pyr	1-Nn-4-(3,4-deH-Pip)
1-3019	4-F-Ph	2-NH ₂ -4-Pyr	1-cPr-4-(3,4-deH-Pip)
1-3020	4-F-Ph	2-NH ₂ -4-Pyr	1-cPn-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3021	4-F-Ph	2-NH ₂ -4-Pyr	1-cHx-4-(3,4-deH-Pip)
1-3022	4-F-Ph	2-NH ₂ -4-Pyr	1-Bn-4-(3,4-deH-Pip)
1-3023	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-4-(3,4-deH-Pip)
1-3024	4-F-Ph	2-NH ₂ -4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
1-3025	4-F-Ph	2-NH ₂ -4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
1-3026	4-F-Ph	2-NH ₂ -4-Pyr	1-Allyl-4-(3,4-deH-Pip)
1-3027	4-F-Ph	2-NH ₂ -4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
1-3028	4-F-Ph	2-NH ₂ -4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
1-3029	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
1-3030	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2,6,6-pentaMe-4-Pip
1-3031	4-F-Ph	2-NH ₂ -4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3032	4-F-Ph	2-NH ₂ -4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3033	4-F-Ph	2-NH ₂ -4-Pyr	7-octaH-Ind
1-3034	4-F-Ph	2-NH ₂ -4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
1-3035	4-F-Ph	2-NH ₂ -4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
1-3036	4-F-Ph	2-NH ₂ -4-Pyr	8-octaH-Qui
1-3037	4-F-Ph	2-NH ₂ -4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
1-3038	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
1-3039	4-F-Ph	2-NH ₂ -4-Pyr	2,2-diMe-4-(4,5-deH-Pip)
1-3040	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)
1-3041	4-F-Ph	2-NH ₂ -4-Pyr	2,6-diMe-4-(3,4-deH-Pip)
1-3042	4-F-Ph	2-NH ₂ -4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
1-3043	4-F-Ph	2-NH ₂ -4-Pyr	2-Me-4-(3,4-deH-Pip)
1-3044	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diMe-4-(3,4-deH-Pip)
1-3045	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
1-3046	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
1-3047	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
1-3048	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3049	4-F-Ph	2-NH ₂ -4-Pyr	2-Et-4-(3,4-deH-Pip)
1-3050	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
1-3051	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
1-3052	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
1-3053	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
1-3054	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
1-3055	4-F-Ph	2-NH ₂ -4-Pyr	2-Pr-4-(3,4-deH-Pip)
1-3056	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
1-3057	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
1-3058	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
1-3059	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
1-3060	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
1-3061	4-F-Ph	2-NH ₂ -4-Pyr	2-Bu-4-(3,4-deH-Pip)
1-3062	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
1-3063	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
1-3064	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
1-3065	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
1-3066	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
1-3067	4-F-Ph	2-NH ₂ -4-Pyr	2-Allyl-4-(3,4-deH-Pip)
1-3068	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
1-3069	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)
1-3070	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)
1-3071	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
1-3072	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
1-3073	4-F-Ph	2-NH ₂ -4-Pyr	2-Bn-4-(3,4-deH-Pip)
1-3074	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)
1-3075	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
1-3076	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
1-3077	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
1-3078	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3079	4-F-Ph	2-NH ₂ -4-Pyr	2-Phet-4-(3,4-deH-Pip)
1-3080	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
1-3081	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
1-3082	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
1-3083	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
1-3084	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
1-3085	4-F-Ph	2-NH ₂ -4-Pyr	2-Me-4-(4,5-deH-Pip)
1-3086	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
1-3087	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
1-3088	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
1-3089	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
1-3090	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
1-3091	4-F-Ph	2-NH ₂ -4-Pyr	2-Et-4-(4,5-deH-Pip)
1-3092	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
1-3093	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
1-3094	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
1-3095	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
1-3096	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
1-3097	4-F-Ph	2-NH ₂ -4-Pyr	2-Pr-4-(4,5-deH-Pip)
1-3098	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
1-3099	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)
1-3100	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPr-4-(4,5-deH-Pip)
1-3101	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
1-3102	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
1-3103	4-F-Ph	2-NH ₂ -4-Pyr	2-Bu-4-(4,5-deH-Pip)
1-3104	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)
1-3105	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
1-3106	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
1-3107	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
1-3108	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3109	4-F-Ph	2-NH ₂ -4-Pyr	2-Allyl-4-(4,5-deH-Pip)
1-3110	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Allyl-4-(4,5-deH-Pip)
1-3111	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
1-3112	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
1-3113	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
1-3114	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
1-3115	4-F-Ph	2-NH ₂ -4-Pyr	2-Bn-4-(4,5-deH-Pip)
1-3116	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
1-3117	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
1-3118	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
1-3119	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
1-3120	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
1-3121	4-F-Ph	2-NH ₂ -4-Pyr	2-Phet-4-(4,5-deH-Pip)
1-3122	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
1-3123	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
1-3124	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
1-3125	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
1-3126	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
1-3127	4-F-Ph	2-MeNH-4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
1-3128	4-F-Ph	2-MeNH-4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
1-3129	4-F-Ph	2-MeNH-4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
1-3130	4-F-Ph	2-MeNH-4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
1-3131	4-F-Ph	2-MeNH-4-Pyr	3-(3,4-deH-Pip)
1-3132	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-(3,4-deH-Pip)
1-3133	4-F-Ph	2-MeNH-4-Pyr	1-Et-4-(3,4-deH-Pip)
1-3134	4-F-Ph	2-MeNH-4-Pyr	1-Pr-4-(3,4-deH-Pip)
1-3135	4-F-Ph	2-MeNH-4-Pyr	1-Pr-4-Pip
1-3136	4-F-Ph	2-MeNH-4-Pyr	1-iPr-4-(3,4-deH-Pip)
1-3137	4-F-Ph	2-MeNH-4-Pyr	1-iPr-4-Pip
1-3138	4-F-Ph	2-MeNH-4-Pyr	1-Bu-4-(3,4-deH-Pip)

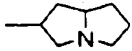

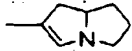
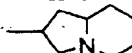
Compound No.	R ¹	R ²	R ³
1-3139	4-F-Ph	2-MeNH-4-Pyr	1-tBu-4-(3,4-deH-Pip)
1-3140	4-F-Ph	2-MeNH-4-Pyr	1-Pn-4-(3,4-deH-Pip)
1-3141	4-F-Ph	2-MeNH-4-Pyr	1-Hx-4-(3,4-deH-Pip)
1-3142	4-F-Ph	2-MeNH-4-Pyr	1-Hp-4-(3,4-deH-Pip)
1-3143	4-F-Ph	2-MeNH-4-Pyr	1-Oc-4-(3,4-deH-Pip)
1-3144	4-F-Ph	2-MeNH-4-Pyr	1-Nn-4-(3,4-deH-Pip)
1-3145	4-F-Ph	2-MeNH-4-Pyr	1-cPr-4-(3,4-deH-Pip)
1-3146	4-F-Ph	2-MeNH-4-Pyr	1-cPn-4-(3,4-deH-Pip)
1-3147	4-F-Ph	2-MeNH-4-Pyr	1-cHx-4-(3,4-deH-Pip)
1-3148	4-F-Ph	2-MeNH-4-Pyr	1-Bn-4-(3,4-deH-Pip)
1-3149	4-F-Ph	2-MeNH-4-Pyr	1-Phet-4-(3,4-deH-Pip)
1-3150	4-F-Ph	2-MeNH-4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
1-3151	4-F-Ph	2-MeNH-4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
1-3152	4-F-Ph	2-MeNH-4-Pyr	1-Allyl-4-(3,4-deH-Pip)
1-3153	4-F-Ph	2-MeNH-4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
1-3154	4-F-Ph	2-MeNH-4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
1-3155	4-F-Ph	2-MeNH-4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
1-3156	4-F-Ph	2-MeNH-4-Pyr	1,2,2,6,6-pentaMe-4-Pip
1-3157	4-F-Ph	2-MeNH-4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3158	4-F-Ph	2-MeNH-4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3159	4-F-Ph	2-MeNH-4-Pyr	7-octaH-Ind
1-3160	4-F-Ph	2-MeNH-4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
1-3161	4-F-Ph	2-MeNH-4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
1-3162	4-F-Ph	2-MeNH-4-Pyr	8-octaH-Qui
1-3163	4-F-Ph	2-MeNH-4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
1-3164	4-F-Ph	2-MeNH-4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
1-3165	4-F-Ph	2-MeNH-4-Pyr	2,2-diMe-4-(4,5-deH-Pip)
1-3166	4-F-Ph	2-MeNH-4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3167	4-F-Ph	2-MeNH-4-Pyr	2,6-diMe-4-(3,4-deH-Pip)
1-3168	4-F-Ph	2-MeNH-4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
1-3169	4-F-Ph	2-MeNH-4-Pyr	2-Me-4-(3,4-deH-Pip)
1-3170	4-F-Ph	2-MeNH-4-Pyr	1,2-diMe-4-(3,4-deH-Pip)
1-3171	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
1-3172	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
1-3173	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
1-3174	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)
1-3175	4-F-Ph	2-MeNH-4-Pyr	2-Et-4-(3,4-deH-Pip)
1-3176	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
1-3177	4-F-Ph	2-MeNH-4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
1-3178	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
1-3179	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
1-3180	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
1-3181	4-F-Ph	2-MeNH-4-Pyr	2-Pr-4-(3,4-deH-Pip)
1-3182	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
1-3183	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
1-3184	4-F-Ph	2-MeNH-4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
1-3185	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
1-3186	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
1-3187	4-F-Ph	2-MeNH-4-Pyr	2-Bu-4-(3,4-deH-Pip)
1-3188	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
1-3189	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
1-3190	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
1-3191	4-F-Ph	2-MeNH-4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
1-3192	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
1-3193	4-F-Ph	2-MeNH-4-Pyr	2-Allyl-4-(3,4-deH-Pip)
1-3194	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
1-3195	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)
1-3196	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3197	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
1-3198	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
1-3199	4-F-Ph	2-MeNH-4-Pyr	2-Bn-4-(3,4-deH-Pip)
1-3200	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)
1-3201	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
1-3202	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
1-3203	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
1-3204	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)
1-3205	4-F-Ph	2-MeNH-4-Pyr	2-Phet-4-(3,4-deH-Pip)
1-3206	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
1-3207	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
1-3208	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
1-3209	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
1-3210	4-F-Ph	2-MeNH-4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
1-3211	4-F-Ph	2-MeNH-4-Pyr	2-Me-4-(4,5-deH-Pip)
1-3212	4-F-Ph	2-MeNH-4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
1-3213	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
1-3214	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
1-3215	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
1-3216	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
1-3217	4-F-Ph	2-MeNH-4-Pyr	2-Et-4-(4,5-deH-Pip)
1-3218	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
1-3219	4-F-Ph	2-MeNH-4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
1-3220	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
1-3221	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
1-3222	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
1-3223	4-F-Ph	2-MeNH-4-Pyr	2-Pr-4-(4,5-deH-Pip)
1-3224	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
1-3225	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)
1-3226	4-F-Ph	2-MeNH-4-Pyr	1,2-diPr-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
1-3227	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
1-3228	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
1-3229	4-F-Ph	2-MeNH-4-Pyr	2-Bu-4-(4,5-deH-Pip)
1-3230	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)
1-3231	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
1-3232	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
1-3233	4-F-Ph	2-MeNH-4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
1-3234	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)
1-3235	4-F-Ph	2-MeNH-4-Pyr	2-Allyl-4-(4,5-deH-Pip)
1-3236	4-F-Ph	2-MeNH-4-Pyr	1-Me-Allyl-4-(4,5-deH-Pip)
1-3237	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
1-3238	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
1-3239	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
1-3240	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
1-3241	4-F-Ph	2-MeNH-4-Pyr	2-Bn-4-(4,5-deH-Pip)
1-3242	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
1-3243	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
1-3244	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
1-3245	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
1-3246	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
1-3247	4-F-Ph	2-MeNH-4-Pyr	2-Phet-4-(4,5-deH-Pip)
1-3248	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
1-3249	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
1-3250	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
1-3251	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
1-3252	4-F-Ph	2-MeNH-4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
1-3253	Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3254	Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3255	3-F-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3256	3-F-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)

Compound No.	R ¹	R ²	R ³
1-3257	3-Cl-Ph	4-PyT	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3258	3-Cl-Ph	4-PyT	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3259	3,4-diF-Ph	4-PyT	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3260	3,4-diF-Ph	4-PyT	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3261	3-CF ₃ -Ph	4-PyT	7-(1,2,3,5,6,8a-hexaH-Ind)
1-3262	3-CF ₃ -Ph	4-PyT	7-(1,2,3,5,8,8a-hexaH-Ind)
1-3263	4-F-Ph	4-PyT	2-Pyrd-CH=CH-
1-3264	4-F-Ph	4-PyT	1-Me-2-Pyrd-CH=CH-
1-3265	4-F-Ph	4-PyT	1-Et-2-Pyrd-CH=CH-
1-3266	4-F-Ph	4-PyT	1-Pr-2-Pyrd-CH=CH-
1-3267	4-F-Ph	4-PyT	2-Pip-CH=CH-
1-3268	4-F-Ph	4-PyT	1-Me-2-Pip-CH=CH-
1-3269	4-F-Ph	4-PyT	1-Et-2-Pip-CH=CH-
1-3270	4-F-Ph	4-PyT	1-Pr-2-Pip-CH=CH-
1-3271	Ph	4-PyT	2-Pyrd-CH=CH-
1-3272	Ph	4-PyT	1-Me-2-Pyrd-CH=CH-
1-3273	Ph	4-PyT	1-Et-2-Pyrd-CH=CH-
1-3274	Ph	4-PyT	1-Pr-2-Pyrd-CH=CH-
1-3275	Ph	4-PyT	2-Pip-CH=CH-
1-3276	Ph	4-PyT	1-Me-2-Pip-CH=CH-
1-3277	Ph	4-PyT	1-Et-2-Pip-CH=CH-
1-3278	Ph	4-PyT	1-Pr-2-Pip-CH=CH-
1-3279	3-F-Ph	4-PyT	2-Pyrd-CH=CH-
1-3280	3-F-Ph	4-PyT	1-Me-2-Pyrd-CH=CH-
1-3281	3-F-Ph	4-PyT	1-Et-2-Pyrd-CH=CH-
1-3282	3-F-Ph	4-PyT	1-Pr-2-Pyrd-CH=CH-
1-3283	3-F-Ph	4-PyT	2-Pip-CH=CH-
1-3284	3-F-Ph	4-PyT	1-Me-2-Pip-CH=CH-
1-3285	3-F-Ph	4-PyT	1-Et-2-Pip-CH=CH-
1-3286	3-F-Ph	4-PyT	1-Pr-2-Pip-CH=CH-

Compound No.	R ¹	R ²	R ³
1-3287	3-Cl-Ph	4-Pyr	2-Pyrd-CH=CH-
1-3288	3-Cl-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
1-3289	3-Cl-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
1-3290	3-Cl-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
1-3291	3-Cl-Ph	4-Pyr	2-Pip-CH=CH-
1-3292	3-Cl-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
1-3293	3-Cl-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
1-3294	3-Cl-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
1-3295	3,4-diF-Ph	4-Pyr	2-Pyrd-CH=CH-
1-3296	3,4-diF-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
1-3297	3,4-diF-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
1-3298	3,4-diF-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
1-3299	3,4-diF-Ph	4-Pyr	2-Pip-CH=CH-
1-3300	3,4-diF-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
1-3301	3,4-diF-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
1-3302	3,4-diF-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
1-3303	3-CF ₃ -Ph	4-Pyr	2-Pyrd-CH=CH-
1-3304	3-CF ₃ -Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
1-3305	3-CF ₃ -Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
1-3306	3-CF ₃ -Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
1-3307	3-CF ₃ -Ph	4-Pyr	2-Pip-CH=CH-
1-3308	3-CF ₃ -Ph	4-Pyr	1-Me-2-Pip-CH=CH-
1-3309	3-CF ₃ -Ph	4-Pyr	1-Et-2-Pip-CH=CH-
1-3310	3-CF ₃ -Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
1-3311	4-F-Ph	4-Pyr	
1-3312	4-F-Ph	4-Pyr	
1-3313	4-F-Ph	4-Pyr	
1-3314	4-F-Ph	4-Pyr	

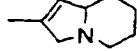
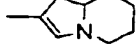
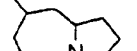
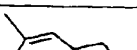
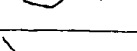
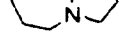
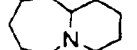
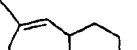
Compound No.	R ¹	R ²	R ³
1-3315	4-F-Ph	4-Pyr	
1-3316	4-F-Ph	4-Pyr	
1-3317	4-F-Ph	4-Pyr	
1-3318	4-F-Ph	4-Pyr	
1-3319	4-F-Ph	4-Pyr	
1-3320	4-F-Ph	4-Pyr	
1-3321	4-F-Ph	4-Pyr	
1-3322	4-F-Ph	4-Pyr	

Table 2

Compound No.	R ¹	R ²	R ³
2-1	Ph	4-Pyr	H ₂ N-CH ₂
2-2	Ph	4-Pyr	H ₂ N-(CH ₂) ₂
2-3	Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-4	Ph	4-Pyr	H ₂ N-(CH ₂) ₄
2-5	Ph	4-Pyr	MeNH-CH ₂
2-6	Ph	4-Pyr	MeNH-(CH ₂) ₂
2-7	Ph	4-Pyr	MeNH-(CH ₂) ₃
2-8	Ph	4-Pyr	MeNH-(CH ₂) ₄
2-9	Ph	4-Pyr	EtNH-CH ₂
2-10	Ph	4-Pyr	EtNH-(CH ₂) ₂
2-11	Ph	4-Pyr	EtNH-(CH ₂) ₃
2-12	Ph	4-Pyr	EtNH-(CH ₂) ₄
2-13	Ph	4-Pyr	Me ₂ N-CH ₂
2-14	Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
2-15	Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-16	Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
2-17	Ph	4-Pyr	1-Azt-CH ₂
2-18	Ph	4-Pyr	1-Azt-(CH ₂) ₂
2-19	Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-20	Ph	4-Pyr	1-Azt-(CH ₂) ₄
2-21	Ph	4-Pyr	1-Pyrd-CH ₂
2-22	Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
2-23	Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-24	Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
2-25	Ph	4-Pyr	1-Pip-CH ₂
2-26	Ph	4-Pyr	1-Pip-(CH ₂) ₂
2-27	Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-28	Ph	4-Pyr	1-Pip-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
2-29	Ph	4-Pyr	1-Mor-CH ₂
2-30	Ph	4-Pyr	1-Mor-(CH ₂) ₂
2-31	Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-32	Ph	4-Pyr	1-Mor-(CH ₂) ₄
2-33	Ph	4-Pyr	1-Tmor-CH ₂
2-34	Ph	4-Pyr	1-Tmor-(CH ₂) ₂
2-35	Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-36	Ph	4-Pyr	1-Tmor-(CH ₂) ₄
2-37	Ph	4-Pyr	1-Piz-CH ₂
2-38	Ph	4-Pyr	1-Piz-(CH ₂) ₂
2-39	Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-40	Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-41	Ph	4-Pyr	1-Piz-(CH ₂) ₄
2-42	Ph	4-Pyr	3-Azt
2-43	Ph	4-Pyr	1-Me-3-Azt
2-44	Ph	4-Pyr	1-Bn-3-Azt
2-45	Ph	4-Pyr	3-Pyrd
2-46	Ph	4-Pyr	1-Me-3-Pyrd
2-47	Ph	4-Pyr	3-Pip
2-48	Ph	4-Pyr	4-Pip
2-49	Ph	4-Pyr	4-(3,4-deH-Pip)
2-50	Ph	4-Pyr	1-Me-4-Pip
2-51	Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-52	Ph	4-Pyr	1-Et-4-Pip
2-53	Ph	4-Pyr	1-Bn-4-Pip
2-54	Ph	4-Pyr	3-Hpip
2-55	Ph	4-Pyr	1-Me-3-Hpip
2-56	Ph	4-Pyr	4-Hpip
2-57	Ph	4-Pyr	1-Me-4-Hpip
2-58	Ph	4-Pyr	2-Mor

Compound No.	R ¹	R ²	R ³
2-59	Ph	4-PyT	1-Me-2-Mor
2-60	Ph	4-PyT	2-Tmor
2-61	Ph	4-PyT	1-Me-2-Tmor
2-62	Ph	4-PyT	1-Piz
2-63	Ph	4-PyT	4-Me-1-Piz
2-64	Ph	4-PyT	2-Piz
2-65	Ph	4-PyT	4-PyT
2-66	Ph	4-PyT	3-PyT
2-67	Ph	4-PyT	2-PyT
2-68	Ph	4-PyT	4-Pym
2-69	Ph	4-PyT	5-Pym
2-70	Ph	4-PyT	2-Pym
2-71	Ph	4-PyT	3-Azt-CH ₂
2-72	Ph	4-PyT	1-Me-3-Azt-CH ₂
2-73	Ph	4-PyT	3-Pyrd-CH ₂
2-74	Ph	4-PyT	1-Me-3-Pyrd-CH ₂
2-75	Ph	4-PyT	4-Pip-CH ₂
2-76	Ph	4-PyT	1-Me-4-Pip-CH ₂
2-77	Ph	4-PyT	3-Hpip-CH ₂
2-78	Ph	4-PyT	1-Me-3-Hpip-CH ₂
2-79	Ph	4-PyT	4-Hpip-CH ₂
2-80	Ph	4-PyT	1-Me-4-Hpip-CH ₂
2-81	Ph	4-PyT	2-Mor-CH ₂
2-82	Ph	4-PyT	1-Me-2-Mor-CH ₂
2-83	Ph	4-PyT	2-Tmor-CH ₂
2-84	Ph	4-PyT	1-Me-2-Tmor-CH ₂
2-85	Ph	4-PyT	1-Piz-CH ₂
2-86	Ph	4-PyT	4-Me-1-Piz-CH ₂
2-87	Ph	4-PyT	2-Piz-CH ₂
2-88	Ph	4-PyT	4-PyT-CH ₂

Compound No.	R ¹	R ²	R ³
2-89	Ph	4-Pyr	3-Pyr-CH ₂
2-90	Ph	4-Pyr	2-Pyr-CH ₂
2-91	Ph	4-Pyr	4-Pym-CH ₂
2-92	Ph	4-Pyr	5-Pym-CH ₂
2-93	Ph	4-Pyr	2-Pym-CH ₂
2-94	Ph	4-Pyr	H ₂ N-CH ₂ CH=CH
2-95	4-F-Ph	4-Pyr	H ₂ N-CH ₂
2-96	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
2-97	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-98	4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
2-99	4-F-Ph	4-Pyr	MeNH-CH ₂
2-100	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₂
2-101	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-102	4-F-Ph	4-Pyr	MeNH-(CH ₂) ₄
2-103	4-F-Ph	4-Pyr	EtNH-CH ₂
2-104	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₂
2-105	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-106	4-F-Ph	4-Pyr	EtNH-(CH ₂) ₄
2-107	4-F-Ph	4-Pyr	Me ₂ N-CH ₂
2-108	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
2-109	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-110	4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
2-111	4-F-Ph	4-Pyr	1-Azt-CH ₂
2-112	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₂
2-113	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-114	4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₄
2-115	4-F-Ph	4-Pyr	1-Pyrd-CH ₂
2-116	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
2-117	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-118	4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
2-119	4-F-Ph	4-Pyr	1-Pip-CH ₂
2-120	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₂
2-121	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-122	4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₄
2-123	4-F-Ph	4-Pyr	1-Mor-CH ₂
2-124	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₂
2-125	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-126	4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₄
2-127	4-F-Ph	4-Pyr	1-Tmor-CH ₂
2-128	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
2-129	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-130	4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
2-131	4-F-Ph	4-Pyr	1-Piz-CH ₂
2-132	4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₂
2-133	4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-134	4-F-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-135	4-F-Ph	4-Pyr	4-Bn-1-Piz-(CH ₂) ₃
2-136	4-F-Ph	4-Pyr	3-Azt
2-137	4-F-Ph	4-Pyr	1-Me-3-Azt
2-138	4-F-Ph	4-Pyr	1-Bn-3-Azt
2-139	4-F-Ph	4-Pyr	3-Pyrd
2-140	4-F-Ph	4-Pyr	1-Me-3-Pyrd
2-141	4-F-Ph	4-Pyr	3-Pip
2-142	4-F-Ph	4-Pyr	4-Pip
2-143	4-F-Ph	4-Pyr	4-(3,4-deH-Pip)
2-144	4-F-Ph	4-Pyr	1-Me-4-Pip
2-145	4-F-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-146	4-F-Ph	4-Pyr	1-Et-4-Pip
2-147	4-F-Ph	4-Pyr	1-Bn-4-Pip
2-148	4-F-Ph	4-Pyr	3-Hpip

Compound No.	R ¹	R ²	R ³
2-149	4-F-Ph	4-Pyr	1-Me-3-Hpip
2-150	4-F-Ph	4-Pyr	4-Hpip
2-151	4-F-Ph	4-Pyr	1-Me-4-Hpip
2-152	4-F-Ph	4-Pyr	2-Mor
2-153	4-F-Ph	4-Pyr	1-Me-2-Mor
2-154	4-F-Ph	4-Pyr	2-Tmor
2-155	4-F-Ph	4-Pyr	1-Me-2-Tmor
2-156	4-F-Ph	4-Pyr	1-Piz
2-157	4-F-Ph	4-Pyr	4-Me-1-Piz
2-158	4-F-Ph	4-Pyr	2-Piz
2-159	4-F-Ph	4-Pyr	4-Pyr
2-160	4-F-Ph	4-Pyr	3-Pyr
2-161	4-F-Ph	4-Pyr	2-Pyr
2-162	4-F-Ph	4-Pyr	4-Pym
2-163	4-F-Ph	4-Pyr	5-Pym
2-164	4-F-Ph	4-Pyr	2-Pym
2-165	4-F-Ph	4-Pyr	3-Azt-CH ₂
2-166	4-F-Ph	4-Pyr	1-Me-3-Azt-CH ₂
2-167	4-F-Ph	4-Pyr	3-Pyrd-CH ₂
2-168	4-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-169	4-F-Ph	4-Pyr	4-Pip-CH ₂
2-170	4-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-171	4-F-Ph	4-Pyr	3-Hpip-CH ₂
2-172	4-F-Ph	4-Pyr	1-Me-3-Hpip-CH ₂
2-173	4-F-Ph	4-Pyr	4-Hpip-CH ₂
2-174	4-F-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
2-175	4-F-Ph	4-Pyr	2-Mor-CH ₂
2-176	4-F-Ph	4-Pyr	1-Me-2-Mor-CH ₂
2-177	4-F-Ph	4-Pyr	2-Tmor-CH ₂
2-178	4-F-Ph	4-Pyr	1-Me-2-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
2-179	4-F-Ph	4-Pyr	1-Piz-CH ₂
2-180	4-F-Ph	4-Pyr	4-Me-1-Piz-CH ₂
2-181	4-F-Ph	4-Pyr	2-Piz-CH ₂
2-182	4-F-Ph	4-Pyr	4-Pyr-CH ₂
2-183	4-F-Ph	4-Pyr	3-Pyr-CH ₂
2-184	4-F-Ph	4-Pyr	2-Pyr-CH ₂
2-185	4-F-Ph	4-Pyr	4-Pym-CH ₂
2-186	4-F-Ph	4-Pyr	5-Pym-CH ₂
2-187	4-F-Ph	4-Pyr	2-Pym-CH ₂
2-188	3-F-Ph	4-Pyr	H ₂ N-CH ₂
2-189	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
2-190	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-191	3-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
2-192	3-F-Ph	4-Pyr	MeNH-CH ₂
2-193	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₂
2-194	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-195	3-F-Ph	4-Pyr	MeNH-(CH ₂) ₄
2-196	3-F-Ph	4-Pyr	EtNH-CH ₂
2-197	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₂
2-198	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-199	3-F-Ph	4-Pyr	EtNH-(CH ₂) ₄
2-200	3-F-Ph	4-Pyr	Me ₂ N-CH ₂
2-201	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
2-202	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-203	3-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
2-204	3-F-Ph	4-Pyr	1-Azt-CH ₂
2-205	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₂
2-206	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-207	3-F-Ph	4-Pyr	1-Azt-(CH ₂) ₄
2-208	3-F-Ph	4-Pyr	1-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-209	3-F-Ph	4-PyT	1-Pyrd-(CH ₂) ₂
2-210	3-F-Ph	4-PyT	1-Pyrd-(CH ₂) ₃
2-211	3-F-Ph	4-PyT	1-Pyrd-(CH ₂) ₄
2-212	3-F-Ph	4-PyT	1-Pip-CH ₂
2-213	3-F-Ph	4-PyT	1-Pip-(CH ₂) ₂
2-214	3-F-Ph	4-PyT	1-Pip-(CH ₂) ₃
2-215	3-F-Ph	4-PyT	1-Pip-(CH ₂) ₄
2-216	3-F-Ph	4-PyT	1-Mor-CH ₂
2-217	3-F-Ph	4-PyT	1-Mor-(CH ₂) ₂
2-218	3-F-Ph	4-PyT	1-Mor-(CH ₂) ₃
2-219	3-F-Ph	4-PyT	1-Mor-(CH ₂) ₄
2-220	3-F-Ph	4-PyT	1-Tmor-CH ₂
2-221	3-F-Ph	4-PyT	1-Tmor-(CH ₂) ₂
2-222	3-F-Ph	4-PyT	1-Tmor-(CH ₂) ₃
2-223	3-F-Ph	4-PyT	1-Tmor-(CH ₂) ₄
2-224	3-F-Ph	4-PyT	1-Piz-CH ₂
2-225	3-F-Ph	4-PyT	1-Piz-(CH ₂) ₂
2-226	3-F-Ph	4-PyT	1-Piz-(CH ₂) ₃
2-227	3-F-Ph	4-PyT	4-Me-1-Piz-(CH ₂) ₃
2-228	3-F-Ph	4-PyT	1-Piz-(CH ₂) ₄
2-229	3-F-Ph	4-PyT	3-Azt
2-230	3-F-Ph	4-PyT	1-Me-3-Azt
2-231	3-F-Ph	4-PyT	1-Bn-3-Azt
2-232	3-F-Ph	4-PyT	3-Pyrd
2-233	3-F-Ph	4-PyT	1-Me-3-Pyrd
2-234	3-F-Ph	4-PyT	3-Pip
2-235	3-F-Ph	4-PyT	4-Pip
2-236	3-F-Ph	4-PyT	4-(3,4-deH-Pip)
2-237	3-F-Ph	4-PyT	1-Me-4-Pip
2-238	3-F-Ph	4-PyT	1-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-239	3-F-Ph	4-Pyr	1-Et-4-Pip
2-240	3-F-Ph	4-Pyr	1-Bn-4-Pip
2-241	3-F-Ph	4-Pyr	3-Hpip
2-242	3-F-Ph	4-Pyr	1-Me-3-Hpip
2-243	3-F-Ph	4-Pyr	4-Hpip
2-244	3-F-Ph	4-Pyr	1-Me-4-Hpip
2-245	3-F-Ph	4-Pyr	2-Mor
2-246	3-F-Ph	4-Pyr	1-Me-2-Mor
2-247	3-F-Ph	4-Pyr	2-Tmor
2-248	3-F-Ph	4-Pyr	1-Me-2-Tmor
2-249	3-F-Ph	4-Pyr	1-Piz
2-250	3-F-Ph	4-Pyr	4-Me-1-Piz
2-251	3-F-Ph	4-Pyr	2-Piz
2-252	3-F-Ph	4-Pyr	4-Pyr
2-253	3-F-Ph	4-Pyr	3-Pyr
2-254	3-F-Ph	4-Pyr	2-Pyr
2-255	3-F-Ph	4-Pyr	4-Pym
2-256	3-F-Ph	4-Pyr	5-Pym
2-257	3-F-Ph	4-Pyr	2-Pym
2-258	3-F-Ph	4-Pyr	3-Azt-CH ₂
2-259	3-F-Ph	4-Pyr	1-Me-3-Azt-CH ₂
2-260	3-F-Ph	4-Pyr	3-Pyrd-CH ₂
2-261	3-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-262	3-F-Ph	4-Pyr	4-Pip-CH ₂
2-263	3-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-264	3-F-Ph	4-Pyr	3-Hpip-CH ₂
2-265	3-F-Ph	4-Pyr	1-Me-3-Hpip-CH ₂
2-266	3-F-Ph	4-Pyr	4-Hpip-CH ₂
2-267	3-F-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
2-268	3-F-Ph	4-Pyr	2-Mor-CH ₂

Compound No.	R ¹	R ²	R ³
2-269	3-F-Ph	4-PyT	1-Me-2-Mor-CH ₂
2-270	3-F-Ph	4-PyT	2-Tmor-CH ₂
2-271	3-F-Ph	4-PyT	1-Me-2-Tmor-CH ₂
2-272	3-F-Ph	4-PyT	1-Piz-CH ₂
2-273	3-F-Ph	4-PyT	4-Me-1-Piz-CH ₂
2-274	3-F-Ph	4-PyT	2-Piz-CH ₂
2-275	3-F-Ph	4-PyT	4-PyT-CH ₂
2-276	3-F-Ph	4-PyT	3-PyT-CH ₂
2-277	3-F-Ph	4-PyT	2-PyT-CH ₂
2-278	3-F-Ph	4-PyT	4-Pym-CH ₂
2-279	3-F-Ph	4-PyT	5-Pym-CH ₂
2-280	3-F-Ph	4-PyT	2-Pym-CH ₂
2-281	3,4-diF-Ph	4-PyT	H ₂ N-CH ₂
2-282	3,4-diF-Ph	4-PyT	H ₂ N-(CH ₂) ₂
2-283	3,4-diF-Ph	4-PyT	H ₂ N-(CH ₂) ₃
2-284	3,4-diF-Ph	4-PyT	H ₂ N-(CH ₂) ₄
2-285	3,4-diF-Ph	4-PyT	MeNH-CH ₂
2-286	3,4-diF-Ph	4-PyT	MeNH-(CH ₂) ₂
2-287	3,4-diF-Ph	4-PyT	MeNH-(CH ₂) ₃
2-288	3,4-diF-Ph	4-PyT	MeNH-(CH ₂) ₄
2-289	3,4-diF-Ph	4-PyT	EtNH-CH ₂
2-290	3,4-diF-Ph	4-PyT	EtNH-(CH ₂) ₂
2-291	3,4-diF-Ph	4-PyT	EtNH-(CH ₂) ₃
2-292	3,4-diF-Ph	4-PyT	EtNH-(CH ₂) ₄
2-293	3,4-diF-Ph	4-PyT	Me ₂ N-CH ₂
2-294	3,4-diF-Ph	4-PyT	Me ₂ N-(CH ₂) ₂
2-295	3,4-diF-Ph	4-PyT	Me ₂ N-(CH ₂) ₃
2-296	3,4-diF-Ph	4-PyT	Me ₂ N-(CH ₂) ₄
2-297	3,4-diF-Ph	4-PyT	1-Azt-CH ₂
2-298	3,4-diF-Ph	4-PyT	1-Azt-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
2-299	3,4-diF-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-300	3,4-diF-Ph	4-Pyr	1-Azt-(CH ₂) ₄
2-301	3,4-diF-Ph	4-Pyr	1-Pyrd-CH ₂
2-302	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
2-303	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-304	3,4-diF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
2-305	3,4-diF-Ph	4-Pyr	1-Pip-CH ₂
2-306	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₂
2-307	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-308	3,4-diF-Ph	4-Pyr	1-Pip-(CH ₂) ₄
2-309	3,4-diF-Ph	4-Pyr	1-Mor-CH ₂
2-310	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₂
2-311	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-312	3,4-diF-Ph	4-Pyr	1-Mor-(CH ₂) ₄
2-313	3,4-diF-Ph	4-Pyr	1-Tmor-CH ₂
2-314	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
2-315	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-316	3,4-diF-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
2-317	3,4-diF-Ph	4-Pyr	1-Piz-CH ₂
2-318	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₂
2-319	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-320	3,4-diF-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-321	3,4-diF-Ph	4-Pyr	1-Piz-(CH ₂) ₄
2-322	3,4-diF-Ph	4-Pyr	3-Azt
2-323	3,4-diF-Ph	4-Pyr	1-Me-3-Azt
2-324	3,4-diF-Ph	4-Pyr	1-Bn-3-Azt
2-325	3,4-diF-Ph	4-Pyr	3-Pyrd
2-326	3,4-diF-Ph	4-Pyr	1-Me-3-Pyrd
2-327	3,4-diF-Ph	4-Pyr	3-Pip
2-328	3,4-diF-Ph	4-Pyr	4-Pip

Compound No.	R ¹	R ²	R ³
2-329	3,4-diF-Ph	4-Pyr	4-(3,4-deH-Pip)
2-330	3,4-diF-Ph	4-Pyr	1-Me-4-Pip
2-331	3,4-diF-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-332	3,4-diF-Ph	4-Pyr	1-Et-4-Pip
2-333	3,4-diF-Ph	4-Pyr	1-Bn-4-Pip
2-334	3,4-diF-Ph	4-Pyr	3-Hpip
2-335	3,4-diF-Ph	4-Pyr	1-Me-3-Hpip
2-336	3,4-diF-Ph	4-Pyr	4-Hpip
2-337	3,4-diF-Ph	4-Pyr	1-Me-4-Hpip
2-338	3,4-diF-Ph	4-Pyr	2-Mor
2-339	3,4-diF-Ph	4-Pyr	1-Me-2-Mor
2-340	3,4-diF-Ph	4-Pyr	2-Tmor
2-341	3,4-diF-Ph	4-Pyr	1-Me-2-Tmor
2-342	3,4-diF-Ph	4-Pyr	1-Piz
2-343	3,4-diF-Ph	4-Pyr	4-Me-1-Piz
2-344	3,4-diF-Ph	4-Pyr	2-Piz
2-345	3,4-diF-Ph	4-Pyr	4-Pyr
2-346	3,4-diF-Ph	4-Pyr	3-Pyr
2-347	3,4-diF-Ph	4-Pyr	2-Pyr
2-348	3,4-diF-Ph	4-Pyr	4-Pym
2-349	3,4-diF-Ph	4-Pyr	5-Pym
2-350	3,4-diF-Ph	4-Pyr	2-Pym
2-351	3,4-diF-Ph	4-Pyr	3-Azt-CH ₂
2-352	3,4-diF-Ph	4-Pyr	1-Me-3-Azt-CH ₂
2-353	3,4-diF-Ph	4-Pyr	3-Pyrd-CH ₂
2-354	3,4-diF-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-355	3,4-diF-Ph	4-Pyr	4-Pip-CH ₂
2-356	3,4-diF-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-357	3,4-diF-Ph	4-Pyr	3-Hpip-CH ₂
2-358	3,4-diF-Ph	4-Pyr	1-Me-3-Hpip-CH ₂

Compound No.	R ¹	R ²	R ³
2-359	3,4-diF-Ph	4-Pyr	4-Hpip-CH ₂
2-360	3,4-diF-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
2-361	3,4-diF-Ph	4-Pyr	2-Mor-CH ₂
2-362	3,4-diF-Ph	4-Pyr	1-Me-2-Mor-CH ₂
2-363	3,4-diF-Ph	4-Pyr	2-Tmor-CH ₂
2-364	3,4-diF-Ph	4-Pyr	1-Me-2-Tmor-CH ₂
2-365	3,4-diF-Ph	4-Pyr	1-Piz-CH ₂
2-366	3,4-diF-Ph	4-Pyr	4-Me-1-Piz-CH ₂
2-367	3,4-diF-Ph	4-Pyr	2-Piz-CH ₂
2-368	3,4-diF-Ph	4-Pyr	4-Pyr-CH ₂
2-369	3,4-diF-Ph	4-Pyr	3-Pyr-CH ₂
2-370	3,4-diF-Ph	4-Pyr	2-Pyr-CH ₂
2-371	3,4-diF-Ph	4-Pyr	4-Pym-CH ₂
2-372	3,4-diF-Ph	4-Pyr	5-Pym-CH ₂
2-373	3,4-diF-Ph	4-Pyr	2-Pym-CH ₂
2-374	3-Cl-Ph	4-Pyr	H ₂ N-CH ₂
2-375	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₂
2-376	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-377	3-Cl-Ph	4-Pyr	H ₂ N-(CH ₂) ₄
2-378	3-Cl-Ph	4-Pyr	MeNH-CH ₂
2-379	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₂
2-380	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-381	3-Cl-Ph	4-Pyr	MeNH-(CH ₂) ₄
2-382	3-Cl-Ph	4-Pyr	EtNH-CH ₂
2-383	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₂
2-384	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-385	3-Cl-Ph	4-Pyr	EtNH-(CH ₂) ₄
2-386	3-Cl-Ph	4-Pyr	Me ₂ N-CH ₂
2-387	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₂
2-388	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-389	3-Cl-Ph	4-Pyr	Me ₂ N-(CH ₂) ₄
2-390	3-Cl-Ph	4-Pyr	1-Azt-CH ₂
2-391	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₂
2-392	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-393	3-Cl-Ph	4-Pyr	1-Azt-(CH ₂) ₄
2-394	3-Cl-Ph	4-Pyr	1-Pyrd-CH ₂
2-395	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₂
2-396	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-397	3-Cl-Ph	4-Pyr	1-Pyrd-(CH ₂) ₄
2-398	3-Cl-Ph	4-Pyr	1-Pip-CH ₂
2-399	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₂
2-400	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-401	3-Cl-Ph	4-Pyr	1-Pip-(CH ₂) ₄
2-402	3-Cl-Ph	4-Pyr	1-Mor-CH ₂
2-403	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₂
2-404	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-405	3-Cl-Ph	4-Pyr	1-Mor-(CH ₂) ₄
2-406	3-Cl-Ph	4-Pyr	1-Tmor-CH ₂
2-407	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₂
2-408	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-409	3-Cl-Ph	4-Pyr	1-Tmor-(CH ₂) ₄
2-410	3-Cl-Ph	4-Pyr	1-Piz-CH ₂
2-411	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₂
2-412	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-413	3-Cl-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-414	3-Cl-Ph	4-Pyr	1-Piz-(CH ₂) ₄
2-415	3-Cl-Ph	4-Pyr	3-Azt
2-416	3-Cl-Ph	4-Pyr	1-Me-3-Azt
2-417	3-Cl-Ph	4-Pyr	1-Bn-3-Azt
2-418	3-Cl-Ph	4-Pyr	3-Pyrd

Compound No.	R ¹	R ²	R ³
2-419	3-Cl-Ph	4-Pyr	1-Me-3-Pyrd
2-420	3-Cl-Ph	4-Pyr	3-Pip
2-421	3-Cl-Ph	4-Pyr	4-Pip
2-422	3-Cl-Ph	4-Pyr	4-(3,4-deH-Pip)
2-423	3-Cl-Ph	4-Pyr	1-Me-4-Pip
2-424	3-Cl-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-425	3-Cl-Ph	4-Pyr	1-Et-4-Pip
2-426	3-Cl-Ph	4-Pyr	1-Bn-4-Pip
2-427	3-Cl-Ph	4-Pyr	3-Hpip
2-428	3-Cl-Ph	4-Pyr	1-Me-3-Hpip
2-429	3-Cl-Ph	4-Pyr	4-Hpip
2-430	3-Cl-Ph	4-Pyr	1-Me-4-Hpip
2-431	3-Cl-Ph	4-Pyr	2-Mor
2-432	3-Cl-Ph	4-Pyr	1-Me-2-Mor
2-433	3-Cl-Ph	4-Pyr	2-Tmor
2-434	3-Cl-Ph	4-Pyr	1-Me-2-Tmor
2-435	3-Cl-Ph	4-Pyr	1-Piz
2-436	3-Cl-Ph	4-Pyr	4-Me-1-Piz
2-437	3-Cl-Ph	4-Pyr	2-Piz
2-438	3-Cl-Ph	4-Pyr	4-Pyr
2-439	3-Cl-Ph	4-Pyr	3-Pyr
2-440	3-Cl-Ph	4-Pyr	2-Pyr
2-441	3-Cl-Ph	4-Pyr	4-Pym
2-442	3-Cl-Ph	4-Pyr	5-Pym
2-443	3-Cl-Ph	4-Pyr	2-Pym
2-444	3-Cl-Ph	4-Pyr	3-Azt-CH ₂
2-445	3-Cl-Ph	4-Pyr	1-Me-3-Azt-CH ₂
2-446	3-Cl-Ph	4-Pyr	3-Pyrd-CH ₂
2-447	3-Cl-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-448	3-Cl-Ph	4-Pyr	4-Pip-CH ₂

Compound No.	R ¹	R ²	R ³
2-449	3-Cl-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-450	3-Cl-Ph	4-Pyr	3-Hpip-CH ₂
2-451	3-Cl-Ph	4-Pyr	1-Me-3-Hpip-CH ₂
2-452	3-Cl-Ph	4-Pyr	4-Hpip-CH ₂
2-453	3-Cl-Ph	4-Pyr	1-Me-4-Hpip-CH ₂
2-454	3-Cl-Ph	4-Pyr	2-Mor-CH ₂
2-455	3-Cl-Ph	4-Pyr	1-Me-2-Mor-CH ₂
2-456	3-Cl-Ph	4-Pyr	2-Tmor-CH ₂
2-457	3-Cl-Ph	4-Pyr	1-Me-2-Tmor-CH ₂
2-458	3-Cl-Ph	4-Pyr	1-Piz-CH ₂
2-459	3-Cl-Ph	4-Pyr	4-Me-1-Piz-CH ₂
2-460	3-Cl-Ph	4-Pyr	2-Piz-CH ₂
2-461	3-Cl-Ph	4-Pyr	4-Pyr-CH ₂
2-462	3-Cl-Ph	4-Pyr	3-Pyr-CH ₂
2-463	3-Cl-Ph	4-Pyr	2-Pyr-CH ₂
2-464	3-Cl-Ph	4-Pyr	4-Pym-CH ₂
2-465	3-Cl-Ph	4-Pyr	5-Pym-CH ₂
2-466	3-Cl-Ph	4-Pyr	2-Pym-CH ₂
2-467	Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
2-468	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
2-469	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-470	Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
2-471	Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
2-472	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
2-473	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-474	Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
2-475	Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
2-476	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
2-477	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-478	Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
2-479	Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
2-480	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
2-481	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-482	Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
2-483	Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
2-484	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
2-485	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-486	Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
2-487	Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
2-488	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
2-489	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-490	Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
2-491	Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
2-492	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
2-493	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-494	Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
2-495	Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
2-496	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
2-497	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-498	Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
2-499	Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
2-500	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
2-501	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-502	Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
2-503	Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-504	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
2-505	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-506	Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-507	Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
2-508	Ph	2-NH ₂ -4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
2-509	Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-510	Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
2-511	Ph	2-NH ₂ -4-Pym	3-Pyrd
2-512	Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-513	Ph	2-NH ₂ -4-Pym	3-Pip
2-514	Ph	2-NH ₂ -4-Pym	4-Pip
2-515	Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-516	Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-517	Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-518	Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
2-519	Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
2-520	Ph	2-NH ₂ -4-Pym	3-Hpip
2-521	Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
2-522	Ph	2-NH ₂ -4-Pym	4-Hpip
2-523	Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
2-524	Ph	2-NH ₂ -4-Pym	2-Mor
2-525	Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
2-526	Ph	2-NH ₂ -4-Pym	2-Tmor
2-527	Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
2-528	Ph	2-NH ₂ -4-Pym	1-Piz
2-529	Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-530	Ph	2-NH ₂ -4-Pym	2-Piz
2-531	Ph	2-NH ₂ -4-Pym	4-Pyr
2-532	Ph	2-NH ₂ -4-Pym	3-Pyr
2-533	Ph	2-NH ₂ -4-Pym	2-Pyr
2-534	Ph	2-NH ₂ -4-Pym	4-Pym
2-535	Ph	2-NH ₂ -4-Pym	5-Pym
2-536	Ph	2-NH ₂ -4-Pym	2-Pym
2-537	Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
2-538	Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂

Compound No.	R ¹	R ²	R ³
2-539	Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-540	Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-541	Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-542	Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-543	Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
2-544	Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
2-545	Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
2-546	Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
2-547	Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
2-548	Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
2-549	Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
2-550	Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
2-551	Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-552	Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
2-553	Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-554	Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-555	Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-556	Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
2-557	Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-558	Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-559	Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-560	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
2-561	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
2-562	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-563	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
2-564	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
2-565	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
2-566	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-567	4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
2-568	4-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂

Compound No.	R ¹	R ²	R ³
2-569	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
2-570	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-571	4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
2-572	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
2-573	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
2-574	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-575	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
2-576	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
2-577	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
2-578	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-579	4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
2-580	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
2-581	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
2-582	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-583	4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
2-584	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
2-585	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
2-586	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-587	4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
2-588	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
2-589	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
2-590	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-591	4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
2-592	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
2-593	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
2-594	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-595	4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
2-596	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-597	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
2-598	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-599	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-600	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
2-601	4-F-Ph	2-NH ₂ -4-Pym	3-Azt
2-602	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-603	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
2-604	4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-605	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-606	4-F-Ph	2-NH ₂ -4-Pym	3-Pip
2-607	4-F-Ph	2-NH ₂ -4-Pym	4-Pip
2-608	4-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-609	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-610	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-611	4-F-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
2-612	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
2-613	4-F-Ph	2-NH ₂ -4-Pym	3-Hpip
2-614	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
2-615	4-F-Ph	2-NH ₂ -4-Pym	4-Hpip
2-616	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
2-617	4-F-Ph	2-NH ₂ -4-Pym	2-Mor
2-618	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
2-619	4-F-Ph	2-NH ₂ -4-Pym	2-Tmor
2-620	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
2-621	4-F-Ph	2-NH ₂ -4-Pym	1-Piz
2-622	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-623	4-F-Ph	2-NH ₂ -4-Pym	2-Piz
2-624	4-F-Ph	2-NH ₂ -4-Pym	4-Pyr
2-625	4-F-Ph	2-NH ₂ -4-Pym	3-Pyr
2-626	4-F-Ph	2-NH ₂ -4-Pym	2-Pyr
2-627	4-F-Ph	2-NH ₂ -4-Pym	4-Pym
2-628	4-F-Ph	2-NH ₂ -4-Pym	5-Pym

Compound No.	R ¹	R ²	R ³
2-629	4-F-Ph	2-NH ₂ -4-Pym	2-Pym
2-630	4-F-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
2-631	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
2-632	4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-633	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-634	4-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-635	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-636	4-F-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
2-637	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
2-638	4-F-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
2-639	4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
2-640	4-F-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
2-641	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
2-642	4-F-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
2-643	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
2-644	4-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-645	4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
2-646	4-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-647	4-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-648	4-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-649	4-F-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
2-650	4-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-651	4-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-652	4-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-653	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
2-654	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
2-655	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-656	3-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
2-657	3-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
2-658	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
2-659	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-660	3-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
2-661	3-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
2-662	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
2-663	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-664	3-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
2-665	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
2-666	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
2-667	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-668	3-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
2-669	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
2-670	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
2-671	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-672	3-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
2-673	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
2-674	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
2-675	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-676	3-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
2-677	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
2-678	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
2-679	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-680	3-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
2-681	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
2-682	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
2-683	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-684	3-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
2-685	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
2-686	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
2-687	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-688	3-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
2-689	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-690	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
2-691	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-692	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-693	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
2-694	3-F-Ph	2-NH ₂ -4-Pym	3-Azt
2-695	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-696	3-F-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
2-697	3-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-698	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-699	3-F-Ph	2-NH ₂ -4-Pym	3-Pip
2-700	3-F-Ph	2-NH ₂ -4-Pym	4-Pip
2-701	3-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-702	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-703	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-704	3-F-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
2-705	3-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
2-706	3-F-Ph	2-NH ₂ -4-Pym	3-Hpip
2-707	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
2-708	3-F-Ph	2-NH ₂ -4-Pym	4-Hpip
2-709	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
2-710	3-F-Ph	2-NH ₂ -4-Pym	2-Mor
2-711	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
2-712	3-F-Ph	2-NH ₂ -4-Pym	2-Tmor
2-713	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
2-714	3-F-Ph	2-NH ₂ -4-Pym	1-Piz
2-715	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-716	3-F-Ph	2-NH ₂ -4-Pym	2-Piz
2-717	3-F-Ph	2-NH ₂ -4-Pym	4-Pyr
2-718	3-F-Ph	2-NH ₂ -4-Pym	3-Pyr

Compound No.	R ¹	R ²	R ³
2-719	3-F-Ph	2-NH ₂ -4-Pym	2-Pyr
2-720	3-F-Ph	2-NH ₂ -4-Pym	4-Pym
2-721	3-F-Ph	2-NH ₂ -4-Pym	5-Pym
2-722	3-F-Ph	2-NH ₂ -4-Pym	2-Pym
2-723	3-F-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
2-724	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
2-725	3-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-726	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-727	3-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-728	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-729	3-F-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
2-730	3-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
2-731	3-F-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
2-732	3-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
2-733	3-F-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
2-734	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
2-735	3-F-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
2-736	3-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
2-737	3-F-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-738	3-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
2-739	3-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-740	3-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-741	3-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-742	3-F-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
2-743	3-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-744	3-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-745	3-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-746	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
2-747	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
2-748	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-749	3,4-diF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
2-750	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
2-751	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
2-752	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-753	3,4-diF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
2-754	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
2-755	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
2-756	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-757	3,4-diF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
2-758	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
2-759	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
2-760	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-761	3,4-diF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
2-762	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
2-763	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
2-764	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-765	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
2-766	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
2-767	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
2-768	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-769	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
2-770	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
2-771	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
2-772	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-773	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
2-774	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
2-775	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂
2-776	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-777	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
2-778	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
2-779	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
2-780	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-781	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
2-782	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-783	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
2-784	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-785	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-786	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
2-787	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Azt
2-788	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-789	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
2-790	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-791	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-792	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pip
2-793	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pip
2-794	3,4-diF-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-795	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-796	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-797	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
2-798	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
2-799	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Hpip
2-800	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
2-801	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Hpip
2-802	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
2-803	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Mor
2-804	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
2-805	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Tmor
2-806	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
2-807	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz
2-808	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz

Compound No.	R ¹	R ²	R ³
2-809	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Piz
2-810	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pyr
2-811	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyr
2-812	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pyr
2-813	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pym
2-814	3,4-diF-Ph	2-NH ₂ -4-Pym	5-Pym
2-815	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pym
2-816	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
2-817	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
2-818	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-819	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-820	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-821	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-822	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
2-823	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
2-824	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
2-825	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
2-826	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
2-827	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
2-828	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
2-829	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
2-830	3,4-diF-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-831	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
2-832	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-833	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-834	3,4-diF-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-835	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂
2-836	3,4-diF-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-837	3,4-diF-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-838	3,4-diF-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂

Compound No.	R ¹	R ²	R ³
2-839	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂
2-840	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₂
2-841	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-842	3-Cl-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₄
2-843	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂
2-844	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₂
2-845	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-846	3-Cl-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₄
2-847	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂
2-848	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₂
2-849	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-850	3-Cl-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₄
2-851	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂
2-852	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₂
2-853	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-854	3-Cl-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₄
2-855	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-CH ₂
2-856	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₂
2-857	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-858	3-Cl-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₄
2-859	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-CH ₂
2-860	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₂
2-861	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-862	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₄
2-863	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-CH ₂
2-864	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₂
2-865	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-866	3-Cl-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₄
2-867	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-CH ₂
2-868	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
2-869	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-870	3-Cl-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₄
2-871	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-CH ₂
2-872	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₂
2-873	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-874	3-Cl-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₄
2-875	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-876	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₂
2-877	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-878	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-879	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₄
2-880	3-Cl-Ph	2-NH ₂ -4-Pym	3-Azt
2-881	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-882	3-Cl-Ph	2-NH ₂ -4-Pym	1-Bn-3-Azt
2-883	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-884	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-885	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pip
2-886	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pip
2-887	3-Cl-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-888	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-889	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-890	3-Cl-Ph	2-NH ₂ -4-Pym	1-Et-4-Pip
2-891	3-Cl-Ph	2-NH ₂ -4-Pym	1-Bn-4-Pip
2-892	3-Cl-Ph	2-NH ₂ -4-Pym	3-Hpip
2-893	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip
2-894	3-Cl-Ph	2-NH ₂ -4-Pym	4-Hpip
2-895	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip
2-896	3-Cl-Ph	2-NH ₂ -4-Pym	2-Mor
2-897	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor
2-898	3-Cl-Ph	2-NH ₂ -4-Pym	2-Tmor

Compound No.	R ¹	R ²	R ³
2-899	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor
2-900	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz
2-901	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-902	3-Cl-Ph	2-NH ₂ -4-Pym	2-Piz
2-903	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pyr
2-904	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyr
2-905	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pyr
2-906	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pym
2-907	3-Cl-Ph	2-NH ₂ -4-Pym	5-Pym
2-908	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pym
2-909	3-Cl-Ph	2-NH ₂ -4-Pym	3-Azt-CH ₂
2-910	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt-CH ₂
2-911	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-912	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-913	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-914	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-915	3-Cl-Ph	2-NH ₂ -4-Pym	3-Hpip-CH ₂
2-916	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-3-Hpip-CH ₂
2-917	3-Cl-Ph	2-NH ₂ -4-Pym	4-Hpip-CH ₂
2-918	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-4-Hpip-CH ₂
2-919	3-Cl-Ph	2-NH ₂ -4-Pym	2-Mor-CH ₂
2-920	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Mor-CH ₂
2-921	3-Cl-Ph	2-NH ₂ -4-Pym	2-Tmor-CH ₂
2-922	3-Cl-Ph	2-NH ₂ -4-Pym	1-Me-2-Tmor-CH ₂
2-923	3-Cl-Ph	2-NH ₂ -4-Pym	1-Piz-CH ₂
2-924	3-Cl-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-CH ₂
2-925	3-Cl-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-926	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-927	3-Cl-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-928	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-929	3-Cl-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-930	3-Cl-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-931	3-Cl-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-932	Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
2-933	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
2-934	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-935	Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
2-936	Ph	2-MeNH-4-Pym	MeNH-CH ₂
2-937	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
2-938	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-939	Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
2-940	Ph	2-MeNH-4-Pym	EtNH-CH ₂
2-941	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
2-942	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-943	Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
2-944	Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
2-945	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
2-946	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-947	Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
2-948	Ph	2-MeNH-4-Pym	1-Azt-CH ₂
2-949	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
2-950	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-951	Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
2-952	Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
2-953	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
2-954	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-955	Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
2-956	Ph	2-MeNH-4-Pym	1-Pip-CH ₂
2-957	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
2-958	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-959	Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
2-960	Ph	2-MeNH-4-Pym	1-Mor-CH ₂
2-961	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
2-962	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-963	Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
2-964	Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
2-965	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
2-966	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-967	Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
2-968	Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-969	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
2-970	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-971	Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-972	Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
2-973	Ph	2-MeNH-4-Pym	3-Azt
2-974	Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-975	Ph	2-MeNH-4-Pym	1-Bn-3-Azt
2-976	Ph	2-MeNH-4-Pym	3-Pyrd
2-977	Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-978	Ph	2-MeNH-4-Pym	3-Pip
2-979	Ph	2-MeNH-4-Pym	4-Pip
2-980	Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-981	Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-982	Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-983	Ph	2-MeNH-4-Pym	1-Et-4-Pip
2-984	Ph	2-MeNH-4-Pym	1-Bn-4-Pip
2-985	Ph	2-MeNH-4-Pym	3-Hpip
2-986	Ph	2-MeNH-4-Pym	1-Me-3-Hpip
2-987	Ph	2-MeNH-4-Pym	4-Hpip
2-988	Ph	2-MeNH-4-Pym	1-Me-4-Hpip

Compound No.	R ¹	R ²	R ³
2-989	Ph	2-MeNH-4-Pym	2-Mor
2-990	Ph	2-MeNH-4-Pym	1-Me-2-Mor
2-991	Ph	2-MeNH-4-Pym	2-Tmor
2-992	Ph	2-MeNH-4-Pym	1-Me-2-Tmor
2-993	Ph	2-MeNH-4-Pym	1-Piz
2-994	Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-995	Ph	2-MeNH-4-Pym	2-Piz
2-996	Ph	2-MeNH-4-Pym	4-Pyr
2-997	Ph	2-MeNH-4-Pym	3-Pyr
2-998	Ph	2-MeNH-4-Pym	2-Pyr
2-999	Ph	2-MeNH-4-Pym	4-Pym
2-1000	Ph	2-MeNH-4-Pym	5-Pym
2-1001	Ph	2-MeNH-4-Pym	2-Pym
2-1002	Ph	2-MeNH-4-Pym	3-Azt-CH ₂
2-1003	Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
2-1004	Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1005	Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1006	Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1007	Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1008	Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
2-1009	Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
2-1010	Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
2-1011	Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
2-1012	Ph	2-MeNH-4-Pym	2-Mor-CH ₂
2-1013	Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
2-1014	Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
2-1015	Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
2-1016	Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1017	Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
2-1018	Ph	2-MeNH-4-Pym	2-Piz-CH ₂

Compound No.	R ¹	R ²	R ³
2-1019	Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1020	Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1021	Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
2-1022	Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1023	Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1024	Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1025	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
2-1026	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
2-1027	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1028	4-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
2-1029	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂
2-1030	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
2-1031	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1032	4-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
2-1033	4-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂
2-1034	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
2-1035	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1036	4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
2-1037	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
2-1038	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
2-1039	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1040	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
2-1041	4-F-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
2-1042	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
2-1043	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1044	4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
2-1045	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
2-1046	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
2-1047	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1048	4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄

Compound No.	R ¹	R ²	R ³
2-1049	4-F-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
2-1050	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
2-1051	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1052	4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
2-1053	4-F-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
2-1054	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
2-1055	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1056	4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
2-1057	4-F-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
2-1058	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
2-1059	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1060	4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
2-1061	4-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1062	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
2-1063	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1064	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1065	4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
2-1066	4-F-Ph	2-MeNH-4-Pym	3-Azt
2-1067	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1068	4-F-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
2-1069	4-F-Ph	2-MeNH-4-Pym	3-Pyrd
2-1070	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1071	4-F-Ph	2-MeNH-4-Pym	3-Pip
2-1072	4-F-Ph	2-MeNH-4-Pym	4-Pip
2-1073	4-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1074	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1075	4-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1076	4-F-Ph	2-MeNH-4-Pym	1-Et-4-Pip
2-1077	4-F-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
2-1078	4-F-Ph	2-MeNH-4-Pym	3-Hpip

Compound No.	R ¹	R ²	R ³
2-1079	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
2-1080	4-F-Ph	2-MeNH-4-Pym	4-Hpip
2-1081	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
2-1082	4-F-Ph	2-MeNH-4-Pym	2-Mor
2-1083	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor
2-1084	4-F-Ph	2-MeNH-4-Pym	2-Tmor
2-1085	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
2-1086	4-F-Ph	2-MeNH-4-Pym	1-Piz
2-1087	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1088	4-F-Ph	2-MeNH-4-Pym	2-Piz
2-1089	4-F-Ph	2-MeNH-4-Pym	4-Pyr
2-1090	4-F-Ph	2-MeNH-4-Pym	3-Pyr
2-1091	4-F-Ph	2-MeNH-4-Pym	2-Pyr
2-1092	4-F-Ph	2-MeNH-4-Pym	4-Pym
2-1093	4-F-Ph	2-MeNH-4-Pym	5-Pym
2-1094	4-F-Ph	2-MeNH-4-Pym	2-Pym
2-1095	4-F-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
2-1096	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
2-1097	4-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1098	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1099	4-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1100	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1101	4-F-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
2-1102	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
2-1103	4-F-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
2-1104	4-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
2-1105	4-F-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
2-1106	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
2-1107	4-F-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
2-1108	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂

Compound No.	R ¹	R ²	R ³
2-1109	4-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1110	4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
2-1111	4-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1112	4-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1113	4-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1114	4-F-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
2-1115	4-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1116	4-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1117	4-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1118	3-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
2-1119	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
2-1120	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1121	3-F-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
2-1122	3-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂
2-1123	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
2-1124	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1125	3-F-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
2-1126	3-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂
2-1127	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
2-1128	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1129	3-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
2-1130	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
2-1131	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
2-1132	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1133	3-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
2-1134	3-F-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
2-1135	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
2-1136	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1137	3-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
2-1138	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-1139	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
2-1140	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1141	3-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
2-1142	3-F-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
2-1143	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
2-1144	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1145	3-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
2-1146	3-F-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
2-1147	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
2-1148	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1149	3-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
2-1150	3-F-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
2-1151	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
2-1152	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1153	3-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
2-1154	3-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1155	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
2-1156	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1157	3-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1158	3-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
2-1159	3-F-Ph	2-MeNH-4-Pym	3-Azt
2-1160	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1161	3-F-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
2-1162	3-F-Ph	2-MeNH-4-Pym	3-Pyrd
2-1163	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1164	3-F-Ph	2-MeNH-4-Pym	3-Pip
2-1165	3-F-Ph	2-MeNH-4-Pym	4-Pip
2-1166	3-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1167	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1168	3-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-1169	3-F-Ph	2-MeNH-4-Pym	1-Et-4-Pip
2-1170	3-F-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
2-1171	3-F-Ph	2-MeNH-4-Pym	3-Hpip
2-1172	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
2-1173	3-F-Ph	2-MeNH-4-Pym	4-Hpip
2-1174	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
2-1175	3-F-Ph	2-MeNH-4-Pym	2-Mor
2-1176	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor
2-1177	3-F-Ph	2-MeNH-4-Pym	2-Tmor
2-1178	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
2-1179	3-F-Ph	2-MeNH-4-Pym	1-Piz
2-1180	3-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1181	3-F-Ph	2-MeNH-4-Pym	2-Piz
2-1182	3-F-Ph	2-MeNH-4-Pym	4-Pyr
2-1183	3-F-Ph	2-MeNH-4-Pym	3-Pyr
2-1184	3-F-Ph	2-MeNH-4-Pym	2-Pyr
2-1185	3-F-Ph	2-MeNH-4-Pym	4-Pym
2-1186	3-F-Ph	2-MeNH-4-Pym	5-Pym
2-1187	3-F-Ph	2-MeNH-4-Pym	2-Pym
2-1188	3-F-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
2-1189	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
2-1190	3-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1191	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1192	3-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1193	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1194	3-F-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
2-1195	3-F-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
2-1196	3-F-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
2-1197	3-F-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
2-1198	3-F-Ph	2-MeNH-4-Pym	2-Mor-CH ₂

Compound No.	R ¹	R ²	R ³
2-1199	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
2-1200	3-F-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
2-1201	3-F-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
2-1202	3-F-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1203	3-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
2-1204	3-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1205	3-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1206	3-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1207	3-F-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
2-1208	3-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1209	3-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1210	3-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1211	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
2-1212	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
2-1213	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1214	3,4-diF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
2-1215	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-CH ₂
2-1216	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
2-1217	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1218	3,4-diF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
2-1219	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-CH ₂
2-1220	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
2-1221	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1222	3,4-diF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
2-1223	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
2-1224	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
2-1225	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1226	3,4-diF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
2-1227	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
2-1228	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂

Compound No.	R ¹	R ²	R ³
2-1229	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1230	3,4-diF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
2-1231	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
2-1232	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
2-1233	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1234	3,4-diF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
2-1235	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
2-1236	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
2-1237	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1238	3,4-diF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
2-1239	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
2-1240	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
2-1241	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1242	3,4-diF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
2-1243	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
2-1244	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
2-1245	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1246	3,4-diF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
2-1247	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1248	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
2-1249	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1250	3,4-diF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1251	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
2-1252	3,4-diF-Ph	2-MeNH-4-Pym	3-Azt
2-1253	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1254	3,4-diF-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
2-1255	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyrd
2-1256	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1257	3,4-diF-Ph	2-MeNH-4-Pym	3-Pip
2-1258	3,4-diF-Ph	2-MeNH-4-Pym	4-Pip

Compound No.	R ¹	R ²	R ³
2-1259	3,4-diF-Ph	2-McNH-4-Pym	4-(3,4-deH-Pip)
2-1260	3,4-diF-Ph	2-McNH-4-Pym	1-Me-4-Pip
2-1261	3,4-diF-Ph	2-McNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1262	3,4-diF-Ph	2-McNH-4-Pym	1-Et-4-Pip
2-1263	3,4-diF-Ph	2-McNH-4-Pym	1-Bn-4-Pip
2-1264	3,4-diF-Ph	2-McNH-4-Pym	3-Hpip
2-1265	3,4-diF-Ph	2-McNH-4-Pym	1-Me-3-Hpip
2-1266	3,4-diF-Ph	2-McNH-4-Pym	4-Hpip
2-1267	3,4-diF-Ph	2-McNH-4-Pym	1-Me-4-Hpip
2-1268	3,4-diF-Ph	2-McNH-4-Pym	2-Mor
2-1269	3,4-diF-Ph	2-McNH-4-Pym	1-Me-2-Mor
2-1270	3,4-diF-Ph	2-McNH-4-Pym	2-Tmor
2-1271	3,4-diF-Ph	2-McNH-4-Pym	1-Me-2-Tmor
2-1272	3,4-diF-Ph	2-McNH-4-Pym	1-Piz
2-1273	3,4-diF-Ph	2-McNH-4-Pym	4-Me-1-Piz
2-1274	3,4-diF-Ph	2-McNH-4-Pym	2-Piz
2-1275	3,4-diF-Ph	2-McNH-4-Pym	4-Pyr
2-1276	3,4-diF-Ph	2-McNH-4-Pym	3-Pyr
2-1277	3,4-diF-Ph	2-McNH-4-Pym	2-Pyr
2-1278	3,4-diF-Ph	2-McNH-4-Pym	4-Pym
2-1279	3,4-diF-Ph	2-McNH-4-Pym	5-Pym
2-1280	3,4-diF-Ph	2-McNH-4-Pym	2-Pym
2-1281	3,4-diF-Ph	2-McNH-4-Pym	3-Azt-CH ₂
2-1282	3,4-diF-Ph	2-McNH-4-Pym	1-Me-3-Azt-CH ₂
2-1283	3,4-diF-Ph	2-McNH-4-Pym	3-Pyrd-CH ₂
2-1284	3,4-diF-Ph	2-McNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1285	3,4-diF-Ph	2-McNH-4-Pym	4-Pip-CH ₂
2-1286	3,4-diF-Ph	2-McNH-4-Pym	1-Me-4-Pip-CH ₂
2-1287	3,4-diF-Ph	2-McNH-4-Pym	3-Hpip-CH ₂
2-1288	3,4-diF-Ph	2-McNH-4-Pym	1-Me-3-Hpip-CH ₂

Compound No.	R ¹	R ²	R ³
2-1289	3,4-diF-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
2-1290	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
2-1291	3,4-diF-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
2-1292	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
2-1293	3,4-diF-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
2-1294	3,4-diF-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
2-1295	3,4-diF-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1296	3,4-diF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
2-1297	3,4-diF-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1298	3,4-diF-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1299	3,4-diF-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1300	3,4-diF-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
2-1301	3,4-diF-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1302	3,4-diF-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1303	3,4-diF-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1304	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂
2-1305	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₂
2-1306	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1307	3-Cl-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₄
2-1308	3-Cl-Ph	2-MeNH-4-Pym	MeNH-CH ₂
2-1309	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₂
2-1310	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1311	3-Cl-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₄
2-1312	3-Cl-Ph	2-MeNH-4-Pym	EtNH-CH ₂
2-1313	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₂
2-1314	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1315	3-Cl-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₄
2-1316	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂
2-1317	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₂
2-1318	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1319	3-Cl-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₄
2-1320	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-CH ₂
2-1321	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₂
2-1322	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1323	3-Cl-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₄
2-1324	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-CH ₂
2-1325	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₂
2-1326	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1327	3-Cl-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₄
2-1328	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-CH ₂
2-1329	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₂
2-1330	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1331	3-Cl-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₄
2-1332	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-CH ₂
2-1333	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₂
2-1334	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1335	3-Cl-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₄
2-1336	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-CH ₂
2-1337	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₂
2-1338	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1339	3-Cl-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₄
2-1340	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1341	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₂
2-1342	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1343	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1344	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₄
2-1345	3-Cl-Ph	2-MeNH-4-Pym	3-Azt
2-1346	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1347	3-Cl-Ph	2-MeNH-4-Pym	1-Bn-3-Azt
2-1348	3-Cl-Ph	2-MeNH-4-Pym	3-Pyrd

Compound No.	R ¹	R ²	R ³
2-1349	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1350	3-Cl-Ph	2-MeNH-4-Pym	3-Pip
2-1351	3-Cl-Ph	2-MeNH-4-Pym	4-Pip
2-1352	3-Cl-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1353	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1354	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1355	3-Cl-Ph	2-MeNH-4-Pym	1-Et-4-Pip
2-1356	3-Cl-Ph	2-MeNH-4-Pym	1-Bn-4-Pip
2-1357	3-Cl-Ph	2-MeNH-4-Pym	3-Hpip
2-1358	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Hpip
2-1359	3-Cl-Ph	2-MeNH-4-Pym	4-Hpip
2-1360	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Hpip
2-1361	3-Cl-Ph	2-MeNH-4-Pym	2-Mor
2-1362	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Mor
2-1363	3-Cl-Ph	2-MeNH-4-Pym	2-Tmor
2-1364	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Tmor
2-1365	3-Cl-Ph	2-MeNH-4-Pym	1-Piz
2-1366	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1367	3-Cl-Ph	2-MeNH-4-Pym	2-Piz
2-1368	3-Cl-Ph	2-MeNH-4-Pym	4-Pyr
2-1369	3-Cl-Ph	2-MeNH-4-Pym	3-Pyr
2-1370	3-Cl-Ph	2-MeNH-4-Pym	2-Pyr
2-1371	3-Cl-Ph	2-MeNH-4-Pym	4-Pym
2-1372	3-Cl-Ph	2-MeNH-4-Pym	5-Pym
2-1373	3-Cl-Ph	2-MeNH-4-Pym	2-Pym
2-1374	3-Cl-Ph	2-MeNH-4-Pym	3-Azt-CH ₂
2-1375	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Azt-CH ₂
2-1376	3-Cl-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1377	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1378	3-Cl-Ph	2-MeNH-4-Pym	4-Pip-CH ₂

Compound No.	R ¹	R ²	R ³
2-1379	3-Cl-Ph	2-MeNH-4-Pym.	1-Me-4-Pip-CH ₂
2-1380	3-Cl-Ph	2-MeNH-4-Pym	3-Hpip-CH ₂
2-1381	3-Cl-Ph	2-MeNH-4-Pym	1-Me-3-Hpip-CH ₂
2-1382	3-Cl-Ph	2-MeNH-4-Pym	4-Hpip-CH ₂
2-1383	3-Cl-Ph	2-MeNH-4-Pym	1-Me-4-Hpip-CH ₂
2-1384	3-Cl-Ph	2-MeNH-4-Pym	2-Mor-CH ₂
2-1385	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Mor-CH ₂
2-1386	3-Cl-Ph	2-MeNH-4-Pym	2-Tmor-CH ₂
2-1387	3-Cl-Ph	2-MeNH-4-Pym	1-Me-2-Tmor-CH ₂
2-1388	3-Cl-Ph	2-MeNH-4-Pym	1-Piz-CH ₂
2-1389	3-Cl-Ph	2-MeNH-4-Pym	4-Me-1-Piz-CH ₂
2-1390	3-Cl-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1391	3-Cl-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1392	3-Cl-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1393	3-Cl-Ph	2-MeNH-4-Pym	2-Pyr-CH ₂
2-1394	3-Cl-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1395	3-Cl-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1396	3-Cl-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1397	3-Cl-4-F-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-1398	3-Cl-4-F-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-1399	3-Cl-4-F-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-1400	3-Cl-4-F-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-1401	3-Cl-4-F-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-1402	3-Cl-4-F-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-1403	3-Cl-4-F-Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-1404	3-Cl-4-F-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-1405	3-Cl-4-F-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-1406	3-Cl-4-F-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-1407	3-Cl-4-F-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1408	3-Cl-4-F-Ph	4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
2-1409	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Azt
2-1410	3-Cl-4-F-Ph	4-Pyr	3-Pyrd
2-1411	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Pyrd
2-1412	3-Cl-4-F-Ph	4-Pyr	4-Pip
2-1413	3-Cl-4-F-Ph	4-Pyr	4-(3,4-deH-Pip)
2-1414	3-Cl-4-F-Ph	4-Pyr	1-Me-4-Pip
2-1415	3-Cl-4-F-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1416	3-Cl-4-F-Ph	4-Pyr	1-Piz
2-1417	3-Cl-4-F-Ph	4-Pyr	4-Me-1-Piz
2-1418	3-Cl-4-F-Ph	4-Pyr	4-Pyr
2-1419	3-Cl-4-F-Ph	4-Pyr	3-Pyr
2-1420	3-Cl-4-F-Ph	4-Pyr	4-Pym
2-1421	3-Cl-4-F-Ph	4-Pyr	5-Pym
2-1422	3-Cl-4-F-Ph	4-Pyr	3-Pyrd-CH ₂
2-1423	3-Cl-4-F-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-1424	3-Cl-4-F-Ph	4-Pyr	4-Pip-CH ₂
2-1425	3-Cl-4-F-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-1426	3-Cl-4-F-Ph	4-Pyr	2-Piz-CH ₂
2-1427	3-Cl-4-F-Ph	4-Pyr	4-Pyr-CH ₂
2-1428	3-Cl-4-F-Ph	4-Pyr	3-Pyr-CH ₂
2-1429	3-Cl-4-F-Ph	4-Pyr	4-Pym-CH ₂
2-1430	3-Cl-4-F-Ph	4-Pyr	5-Pym-CH ₂
2-1431	3-Cl-4-F-Ph	4-Pyr	2-Pym-CH ₂
2-1432	3,4,5-triF-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-1433	3,4,5-triF-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-1434	3,4,5-triF-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-1435	3,4,5-triF-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-1436	3,4,5-triF-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-1437	3,4,5-triF-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-1438	3,4,5-triF-Ph	4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1439	3,4,5-triF-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-1440	3,4,5-triF-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-1441	3,4,5-triF-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-1442	3,4,5-triF-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1443	3,4,5-triF-Ph	4-Pyr	3-Azt
2-1444	3,4,5-triF-Ph	4-Pyr	1-Me-3-Azt
2-1445	3,4,5-triF-Ph	4-Pyr	3-Pyrd
2-1446	3,4,5-triF-Ph	4-Pyr	1-Me-3-Pyrd
2-1447	3,4,5-triF-Ph	4-Pyr	4-Pip
2-1448	3,4,5-triF-Ph	4-Pyr	4-(3,4-deH-Pip)
2-1449	3,4,5-triF-Ph	4-Pyr	1-Me-4-Pip
2-1450	3,4,5-triF-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1451	3,4,5-triF-Ph	4-Pyr	1-Piz
2-1452	3,4,5-triF-Ph	4-Pyr	4-Me-1-Piz
2-1453	3,4,5-triF-Ph	4-Pyr	4-Pyr
2-1454	3,4,5-triF-Ph	4-Pyr	3-Pyr
2-1455	3,4,5-triF-Ph	4-Pyr	4-Pym
2-1456	3,4,5-triF-Ph	4-Pyr	5-Pym
2-1457	3,4,5-triF-Ph	4-Pyr	3-Pyrd-CH ₂
2-1458	3,4,5-triF-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-1459	3,4,5-triF-Ph	4-Pyr	4-Pip-CH ₂
2-1460	3,4,5-triF-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-1461	3,4,5-triF-Ph	4-Pyr	2-Piz-CH ₂
2-1462	3,4,5-triF-Ph	4-Pyr	4-Pyr-CH ₂
2-1463	3,4,5-triF-Ph	4-Pyr	3-Pyr-CH ₂
2-1464	3,4,5-triF-Ph	4-Pyr	4-Pym-CH ₂
2-1465	3,4,5-triF-Ph	4-Pyr	5-Pym-CH ₂
2-1466	3,4,5-triF-Ph	4-Pyr	2-Pym-CH ₂
2-1467	3-CF ₃ -Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-1468	3-CF ₃ -Ph	4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1469	3-CF ₃ -Ph	4-Pyr	EtNH-(CH ₂) ₃
2-1470	3-CF ₃ -Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-1471	3-CF ₃ -Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-1472	3-CF ₃ -Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-1473	3-CF ₃ -Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-1474	3-CF ₃ -Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-1475	3-CF ₃ -Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-1476	3-CF ₃ -Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-1477	3-CF ₃ -Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1478	3-CF ₃ -Ph	4-Pyr	3-Azt
2-1479	3-CF ₃ -Ph	4-Pyr	1-Me-3-Azt
2-1480	3-CF ₃ -Ph	4-Pyr	3-Pyrd
2-1481	3-CF ₃ -Ph	4-Pyr	1-Me-3-Pyrd
2-1482	3-CF ₃ -Ph	4-Pyr	4-Pip
2-1483	3-CF ₃ -Ph	4-Pyr	4-(3,4-deH-Pip)
2-1484	3-CF ₃ -Ph	4-Pyr	1-Me-4-Pip
2-1485	3-CF ₃ -Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1486	3-CF ₃ -Ph	4-Pyr	1-Piz
2-1487	3-CF ₃ -Ph	4-Pyr	4-Me-1-Piz
2-1488	3-CF ₃ -Ph	4-Pyr	4-Pyr
2-1489	3-CF ₃ -Ph	4-Pyr	3-Pyr
2-1490	3-CF ₃ -Ph	4-Pyr	4-Pym
2-1491	3-CF ₃ -Ph	4-Pyr	5-Pym
2-1492	3-CF ₃ -Ph	4-Pyr	3-Pyrd-CH ₂
2-1493	3-CF ₃ -Ph	4-Pyr	1-Me-3-Pyrd-CH ₂
2-1494	3-CF ₃ -Ph	4-Pyr	4-Pip-CH ₂
2-1495	3-CF ₃ -Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-1496	3-CF ₃ -Ph	4-Pyr	2-Piz-CH ₂
2-1497	3-CF ₃ -Ph	4-Pyr	4-Pyr-CH ₂
2-1498	3-CF ₃ -Ph	4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-1499	3-CF ₃ -Ph	4-Pyr	4-Pym-CH ₂
2-1500	3-CF ₃ -Ph	4-Pyr	5-Pym-CH ₂
2-1501	3-CF ₃ -Ph	4-Pyr	2-Pym-CH ₂
2-1502	3-CHF ₂ O-Ph	4-Pyr	H ₂ N-(CH ₂) ₃
2-1503	3-CHF ₂ O-Ph	4-Pyr	MeNH-(CH ₂) ₃
2-1504	3-CHF ₂ O-Ph	4-Pyr	EtNH-(CH ₂) ₃
2-1505	3-CHF ₂ O-Ph	4-Pyr	Me ₂ N-(CH ₂) ₃
2-1506	3-CHF ₂ O-Ph	4-Pyr	1-Azt-(CH ₂) ₃
2-1507	3-CHF ₂ O-Ph	4-Pyr	1-Pyrd-(CH ₂) ₃
2-1508	3-CHF ₂ O-Ph	4-Pyr	1-Pip-(CH ₂) ₃
2-1509	3-CHF ₂ O-Ph	4-Pyr	1-Mor-(CH ₂) ₃
2-1510	3-CHF ₂ O-Ph	4-Pyr	1-Tmor-(CH ₂) ₃
2-1511	3-CHF ₂ O-Ph	4-Pyr	1-Piz-(CH ₂) ₃
2-1512	3-CHF ₂ O-Ph	4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1513	3-CHF ₂ O-Ph	4-Pyr	3-Azt
2-1514	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Azt
2-1515	3-CHF ₂ O-Ph	4-Pyr	3-Pyrd
2-1516	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Pyrd
2-1517	3-CHF ₂ O-Ph	4-Pyr	4-Pip
2-1518	3-CHF ₂ O-Ph	4-Pyr	4-(3,4-deH-Pip)
2-1519	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-Pip
2-1520	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1521	3-CHF ₂ O-Ph	4-Pyr	1-Piz
2-1522	3-CHF ₂ O-Ph	4-Pyr	4-Me-1-Piz
2-1523	3-CHF ₂ O-Ph	4-Pyr	4-Pyr
2-1524	3-CHF ₂ O-Ph	4-Pyr	3-Pyr
2-1525	3-CHF ₂ O-Ph	4-Pyr	4-Pym
2-1526	3-CHF ₂ O-Ph	4-Pyr	5-Pym
2-1527	3-CHF ₂ O-Ph	4-Pyr	3-Pyrd-CH ₂
2-1528	3-CHF ₂ O-Ph	4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-1529	3-CHF ₂ O-Ph	4-Pyr	4-Pip-CH ₂
2-1530	3-CHF ₂ O-Ph	4-Pyr	1-Me-4-Pip-CH ₂
2-1531	3-CHF ₂ O-Ph	4-Pyr	2-Piz-CH ₂
2-1532	3-CHF ₂ O-Ph	4-Pyr	4-Pyr-CH ₂
2-1533	3-CHF ₂ O-Ph	4-Pyr	3-Pyr-CH ₂
2-1534	3-CHF ₂ O-Ph	4-Pyr	4-Pym-CH ₂
2-1535	3-CHF ₂ O-Ph	4-Pyr	5-Pym-CH ₂
2-1536	3-CHF ₂ O-Ph	4-Pyr	2-Pym-CH ₂
2-1537	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-1538	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-1539	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-1540	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-1541	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-1542	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-1543	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-1544	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-1545	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-1546	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-1547	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1548	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Azt
2-1549	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-1550	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-1551	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-1552	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pip
2-1553	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-1554	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-1555	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-1556	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Piz
2-1557	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-1558	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
2-1559	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyr
2-1560	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pym
2-1561	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	5-Pym
2-1562	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-1563	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-1564	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-1565	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-1566	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-1567	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-1568	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-1569	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-1570	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-1571	3-Cl-4-F-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-1572	3,4,5-triF-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-1573	3,4,5-triF-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-1574	3,4,5-triF-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-1575	3,4,5-triF-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-1576	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-1577	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-1578	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-1579	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-1580	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-1581	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-1582	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1583	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Azt
2-1584	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-1585	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-1586	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-1587	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pip
2-1588	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-1589	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-1590	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-1591	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Piz
2-1592	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-1593	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pyr
2-1594	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyr
2-1595	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pym
2-1596	3,4,5-triF-Ph	2-NH ₂ -4-Pym	5-Pym
2-1597	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-1598	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-1599	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-1600	3,4,5-triF-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-1601	3,4,5-triF-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-1602	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-1603	3,4,5-triF-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-1604	3,4,5-triF-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-1605	3,4,5-triF-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-1606	3,4,5-triF-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-1607	3-CF ₃ -Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-1608	3-CF ₃ -Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-1609	3-CF ₃ -Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-1610	3-CF ₃ -Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-1611	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-1612	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-1613	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃
2-1614	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-1615	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-1616	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-1617	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1618	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
2-1619	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-1620	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyrd
2-1621	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-1622	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pip
2-1623	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-1624	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-1625	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-1626	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Piz
2-1627	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Me-1-Piz
2-1628	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pyr
2-1629	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyr
2-1630	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pym
2-1631	3-CF ₃ -Ph	2-NH ₂ -4-Pym	5-Pym
2-1632	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-1633	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-1634	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-1635	3-CF ₃ -Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-1636	3-CF ₃ -Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-1637	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-1638	3-CF ₃ -Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-1639	3-CF ₃ -Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-1640	3-CF ₃ -Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-1641	3-CF ₃ -Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-1642	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	H ₂ N-(CH ₂) ₃
2-1643	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	MeNH-(CH ₂) ₃
2-1644	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	EtNH-(CH ₂) ₃
2-1645	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	Me ₂ N-(CH ₂) ₃
2-1646	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Azt-(CH ₂) ₃
2-1647	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Pyrd-(CH ₂) ₃
2-1648	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1649	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Mor-(CH ₂) ₃
2-1650	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Tmor-(CH ₂) ₃
2-1651	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Piz-(CH ₂) ₃
2-1652	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1653	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Azt
2-1654	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Azt
2-1655	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyrd
2-1656	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd
2-1657	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pip
2-1658	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-(3,4-deH-Pip)
2-1659	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip
2-1660	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-(3,4-deH-Pip)
2-1661	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Piz
2-1662	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Mc-1-Piz
2-1663	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pyr
2-1664	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyr
2-1665	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pym
2-1666	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	5-Pym
2-1667	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyrd-CH ₂
2-1668	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-3-Pyrd-CH ₂
2-1669	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pip-CH ₂
2-1670	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	1-Me-4-Pip-CH ₂
2-1671	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	2-Piz-CH ₂
2-1672	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pyr-CH ₂
2-1673	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	3-Pyr-CH ₂
2-1674	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	4-Pym-CH ₂
2-1675	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	5-Pym-CH ₂
2-1676	3-CHF ₂ O-Ph	2-NH ₂ -4-Pym	2-Pym-CH ₂
2-1677	3-Cl-4-F-Ph	2-McNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1678	3-Cl-4-F-Ph	2-McNH-4-Pym	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1679	3-Cl-4-F-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1680	3-Cl-4-F-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1681	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1682	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1683	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1684	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1685	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1686	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1687	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1688	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Azt
2-1689	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1690	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyrd
2-1691	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1692	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pip
2-1693	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1694	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1695	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1696	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Piz
2-1697	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1698	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pyr
2-1699	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyr
2-1700	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pym
2-1701	3-Cl-4-F-Ph	2-MeNH-4-Pym	5-Pym
2-1702	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1703	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1704	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1705	3-Cl-4-F-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1706	3-Cl-4-F-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1707	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1708	3-Cl-4-F-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-1709	3-Cl-4-F-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1710	3-Cl-4-F-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1711	3-Cl-4-F-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1712	3,4,5-triF-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1713	3,4,5-triF-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1714	3,4,5-triF-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1715	3,4,5-triF-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1716	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1717	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1718	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1719	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1720	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1721	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1722	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1723	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Azt
2-1724	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1725	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyrd
2-1726	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1727	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pip
2-1728	3,4,5-triF-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1729	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1730	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1731	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Piz
2-1732	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1733	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pyr
2-1734	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyr
2-1735	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pym
2-1736	3,4,5-triF-Ph	2-MeNH-4-Pym	5-Pym
2-1737	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1738	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-1739	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1740	3,4,5-triF-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1741	3,4,5-triF-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1742	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1743	3,4,5-triF-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1744	3,4,5-triF-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1745	3,4,5-triF-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1746	3,4,5-triF-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1747	3-CF ₃ -Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1748	3-CF ₃ -Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1749	3-CF ₃ -Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1750	3-CF ₃ -Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1751	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1752	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1753	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1754	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1755	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1756	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1757	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1758	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Azt
2-1759	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1760	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyrd
2-1761	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1762	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pip
2-1763	3-CF ₃ -Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)
2-1764	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1765	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1766	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Piz
2-1767	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1768	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
2-1769	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyr
2-1770	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pym
2-1771	3-CF ₃ -Ph	2-MeNH-4-Pym	5-Pym
2-1772	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1773	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1774	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1775	3-CF ₃ -Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1776	3-CF ₃ -Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1777	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1778	3-CF ₃ -Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1779	3-CF ₃ -Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1780	3-CF ₃ -Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1781	3-CF ₃ -Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1782	3-CHF ₂ O-Ph	2-MeNH-4-Pym	H ₂ N-(CH ₂) ₃
2-1783	3-CHF ₂ O-Ph	2-MeNH-4-Pym	MeNH-(CH ₂) ₃
2-1784	3-CHF ₂ O-Ph	2-MeNH-4-Pym	EtNH-(CH ₂) ₃
2-1785	3-CHF ₂ O-Ph	2-MeNH-4-Pym	Me ₂ N-(CH ₂) ₃
2-1786	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Azt-(CH ₂) ₃
2-1787	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Pyrd-(CH ₂) ₃
2-1788	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Pip-(CH ₂) ₃
2-1789	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Mor-(CH ₂) ₃
2-1790	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Tmor-(CH ₂) ₃
2-1791	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Piz-(CH ₂) ₃
2-1792	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-1793	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Azt
2-1794	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Azt
2-1795	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyrd
2-1796	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd
2-1797	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pip
2-1798	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-1799	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-Pip
2-1800	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-(3,4-deH-Pip)
2-1801	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Piz
2-1802	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Me-1-Piz
2-1803	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pyr
2-1804	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyr
2-1805	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pym
2-1806	3-CHF ₂ O-Ph	2-MeNH-4-Pym	5-Pym
2-1807	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyrd-CH ₂
2-1808	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-3-Pyrd-CH ₂
2-1809	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pip-CH ₂
2-1810	3-CHF ₂ O-Ph	2-MeNH-4-Pym	1-Me-4-Pip-CH ₂
2-1811	3-CHF ₂ O-Ph	2-MeNH-4-Pym	2-Piz-CH ₂
2-1812	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pyr-CH ₂
2-1813	3-CHF ₂ O-Ph	2-MeNH-4-Pym	3-Pyr-CH ₂
2-1814	3-CHF ₂ O-Ph	2-MeNH-4-Pym	4-Pym-CH ₂
2-1815	3-CHF ₂ O-Ph	2-MeNH-4-Pym	5-Pym-CH ₂
2-1816	3-CHF ₂ O-Ph	2-MeNH-4-Pym	2-Pym-CH ₂
2-1817	Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
2-1818	Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
2-1819	Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
2-1820	Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
2-1821	Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
2-1822	Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
2-1823	Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
2-1824	Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
2-1825	Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
2-1826	Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
2-1827	Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1828	Ph	2-NH ₂ -4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
2-1829	Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
2-1830	Ph	2-NH ₂ -4-Pyr	3-Pyrd
2-1831	Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
2-1832	Ph	2-NH ₂ -4-Pyr	4-Pip
2-1833	Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
2-1834	Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
2-1835	Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1836	Ph	2-NH ₂ -4-Pyr	1-Piz
2-1837	Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
2-1838	Ph	2-NH ₂ -4-Pyr	4-Pyr
2-1839	Ph	2-NH ₂ -4-Pyr	3-Pyr
2-1840	Ph	2-NH ₂ -4-Pyr	4-Pym
2-1841	Ph	2-NH ₂ -4-Pyr	5-Pym
2-1842	Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
2-1843	Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
2-1844	Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
2-1845	Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
2-1846	Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
2-1847	Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
2-1848	Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
2-1849	Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
2-1850	Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
2-1851	Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
2-1852	4-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
2-1853	4-F-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
2-1854	4-F-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
2-1855	4-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
2-1856	4-F-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
2-1857	4-F-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
2-1858	4-F-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1859	4-F-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
2-1860	4-F-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
2-1861	4-F-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
2-1862	4-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1863	4-F-Ph	2-NH ₂ -4-Pyr	3-Azt
2-1864	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
2-1865	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd
2-1866	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
2-1867	4-F-Ph	2-NH ₂ -4-Pyr	4-Pip
2-1868	4-F-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
2-1869	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
2-1870	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1871	4-F-Ph	2-NH ₂ -4-Pyr	1-Piz
2-1872	4-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
2-1873	4-F-Ph	2-NH ₂ -4-Pyr	4-Pyr
2-1874	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyr
2-1875	4-F-Ph	2-NH ₂ -4-Pyr	4-Pym
2-1876	4-F-Ph	2-NH ₂ -4-Pyr	5-Pym
2-1877	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
2-1878	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
2-1879	4-F-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
2-1880	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
2-1881	4-F-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
2-1882	4-F-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
2-1883	4-F-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
2-1884	4-F-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
2-1885	4-F-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
2-1886	4-F-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
2-1887	3-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
2-1888	3-F-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-1889	3-F-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
2-1890	3-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
2-1891	3-F-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
2-1892	3-F-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
2-1893	3-F-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
2-1894	3-F-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
2-1895	3-F-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
2-1896	3-F-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
2-1897	3-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1898	3-F-Ph	2-NH ₂ -4-Pyr	3-Azt
2-1899	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
2-1900	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd
2-1901	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
2-1902	3-F-Ph	2-NH ₂ -4-Pyr	4-Pip
2-1903	3-F-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
2-1904	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
2-1905	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1906	3-F-Ph	2-NH ₂ -4-Pyr	1-Piz
2-1907	3-F-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
2-1908	3-F-Ph	2-NH ₂ -4-Pyr	4-Pyr
2-1909	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyr
2-1910	3-F-Ph	2-NH ₂ -4-Pyr	4-Pym
2-1911	3-F-Ph	2-NH ₂ -4-Pyr	5-Pym
2-1912	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
2-1913	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
2-1914	3-F-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
2-1915	3-F-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
2-1916	3-F-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
2-1917	3-F-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
2-1918	3-F-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-1919	3-F-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
2-1920	3-F-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
2-1921	3-F-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
2-1922	3,4-diF-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
2-1923	3,4-diF-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
2-1924	3,4-diF-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
2-1925	3,4-diF-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
2-1926	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
2-1927	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
2-1928	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
2-1929	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
2-1930	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
2-1931	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
2-1932	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1933	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Azt
2-1934	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
2-1935	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyrd
2-1936	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
2-1937	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pip
2-1938	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
2-1939	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Mc-4-Pip
2-1940	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1941	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Piz
2-1942	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
2-1943	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pyr
2-1944	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyr
2-1945	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pym
2-1946	3,4-diF-Ph	2-NH ₂ -4-Pyr	5-Pym
2-1947	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
2-1948	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-1949	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
2-1950	3,4-diF-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
2-1951	3,4-diF-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
2-1952	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
2-1953	3,4-diF-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
2-1954	3,4-diF-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
2-1955	3,4-diF-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
2-1956	3,4-diF-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
2-1957	3-Cl-Ph	2-NH ₂ -4-Pyr	H ₂ N-(CH ₂) ₃
2-1958	3-Cl-Ph	2-NH ₂ -4-Pyr	MeNH-(CH ₂) ₃
2-1959	3-Cl-Ph	2-NH ₂ -4-Pyr	EtNH-(CH ₂) ₃
2-1960	3-Cl-Ph	2-NH ₂ -4-Pyr	Me ₂ N-(CH ₂) ₃
2-1961	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Azt-(CH ₂) ₃
2-1962	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Pyrd-(CH ₂) ₃
2-1963	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Pip-(CH ₂) ₃
2-1964	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Mor-(CH ₂) ₃
2-1965	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Tmor-(CH ₂) ₃
2-1966	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Piz-(CH ₂) ₃
2-1967	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-1968	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Azt
2-1969	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Azt
2-1970	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyrd
2-1971	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd
2-1972	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pip
2-1973	3-Cl-Ph	2-NH ₂ -4-Pyr	4-(3,4-deH-Pip)
2-1974	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip
2-1975	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-(3,4-deH-Pip)
2-1976	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Piz
2-1977	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Me-1-Piz
2-1978	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pyr

Compound No.	R ¹	R ²	R ³
2-1979	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyr
2-1980	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pym
2-1981	3-Cl-Ph	2-NH ₂ -4-Pyr	5-Pym
2-1982	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyrd-CH ₂
2-1983	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-3-Pyrd-CH ₂
2-1984	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pip-CH ₂
2-1985	3-Cl-Ph	2-NH ₂ -4-Pyr	1-Me-4-Pip-CH ₂
2-1986	3-Cl-Ph	2-NH ₂ -4-Pyr	2-Piz-CH ₂
2-1987	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pyr-CH ₂
2-1988	3-Cl-Ph	2-NH ₂ -4-Pyr	3-Pyr-CH ₂
2-1989	3-Cl-Ph	2-NH ₂ -4-Pyr	4-Pym-CH ₂
2-1990	3-Cl-Ph	2-NH ₂ -4-Pyr	5-Pym-CH ₂
2-1991	3-Cl-Ph	2-NH ₂ -4-Pyr	2-Pym-CH ₂
2-1992	Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
2-1993	Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
2-1994	Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
2-1995	Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
2-1996	Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
2-1997	Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
2-1998	Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
2-1999	Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
2-2000	Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
2-2001	Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
2-2002	Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-2003	Ph	2-MeNH-4-Pyr	3-Azt
2-2004	Ph	2-MeNH-4-Pyr	1-Me-3-Azt
2-2005	Ph	2-MeNH-4-Pyr	3-Pyrd
2-2006	Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
2-2007	Ph	2-MeNH-4-Pyr	4-Pip
2-2008	Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2009	Ph	2-MeNH-4-Pyr	1-Me-4-Pip
2-2010	Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
2-2011	Ph	2-MeNH-4-Pyr	1-Piz
2-2012	Ph	2-MeNH-4-Pyr	4-Me-1-Piz
2-2013	Ph	2-MeNH-4-Pyr	4-Pyr
2-2014	Ph	2-MeNH-4-Pyr	3-Pyr
2-2015	Ph	2-MeNH-4-Pyr	4-Pym
2-2016	Ph	2-MeNH-4-Pyr	5-Pym
2-2017	Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
2-2018	Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
2-2019	Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
2-2020	Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
2-2021	Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
2-2022	Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
2-2023	Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
2-2024	Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
2-2025	Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
2-2026	Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
2-2027	4-F-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
2-2028	4-F-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
2-2029	4-F-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
2-2030	4-F-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
2-2031	4-F-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
2-2032	4-F-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
2-2033	4-F-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
2-2034	4-F-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
2-2035	4-F-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
2-2036	4-F-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
2-2037	4-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-2038	4-F-Ph	2-MeNH-4-Pyr	3-Azt

Compound No.	R ¹	R ²	R ³
2-2039	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
2-2040	4-F-Ph	2-MeNH-4-Pyr	3-Pyrd
2-2041	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
2-2042	4-F-Ph	2-MeNH-4-Pyr	4-Pip
2-2043	4-F-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
2-2044	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
2-2045	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
2-2046	4-F-Ph	2-MeNH-4-Pyr	1-Piz
2-2047	4-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
2-2048	4-F-Ph	2-MeNH-4-Pyr	4-Pyr
2-2049	4-F-Ph	2-MeNH-4-Pyr	3-Pyr
2-2050	4-F-Ph	2-MeNH-4-Pyr	4-Pym
2-2051	4-F-Ph	2-MeNH-4-Pyr	5-Pym
2-2052	4-F-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
2-2053	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
2-2054	4-F-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
2-2055	4-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
2-2056	4-F-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
2-2057	4-F-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
2-2058	4-F-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
2-2059	4-F-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
2-2060	4-F-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
2-2061	4-F-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
2-2062	3-F-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
2-2063	3-F-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
2-2064	3-F-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
2-2065	3-F-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
2-2066	3-F-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
2-2067	3-F-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
2-2068	3-F-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-2069	3-F-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
2-2070	3-F-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
2-2071	3-F-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
2-2072	3-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-2073	3-F-Ph	2-MeNH-4-Pyr	3-Azt
2-2074	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
2-2075	3-F-Ph	2-MeNH-4-Pyr	3-Pyrd
2-2076	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
2-2077	3-F-Ph	2-MeNH-4-Pyr	4-Pip
2-2078	3-F-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
2-2079	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
2-2080	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
2-2081	3-F-Ph	2-MeNH-4-Pyr	1-Piz
2-2082	3-F-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
2-2083	3-F-Ph	2-MeNH-4-Pyr	4-Pyr
2-2084	3-F-Ph	2-MeNH-4-Pyr	3-Pyr
2-2085	3-F-Ph	2-MeNH-4-Pyr	4-Pym
2-2086	3-F-Ph	2-MeNH-4-Pyr	5-Pym
2-2087	3-F-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
2-2088	3-F-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
2-2089	3-F-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
2-2090	3-F-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
2-2091	3-F-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
2-2092	3-F-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
2-2093	3-F-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
2-2094	3-F-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
2-2095	3-F-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
2-2096	3-F-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
2-2097	3,4-diF-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
2-2098	3,4-diF-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-2099	3,4-diF-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
2-2100	3,4-diF-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
2-2101	3,4-diF-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
2-2102	3,4-diF-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
2-2103	3,4-diF-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
2-2104	3,4-diF-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
2-2105	3,4-diF-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
2-2106	3,4-diF-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
2-2107	3,4-diF-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-2108	3,4-diF-Ph	2-MeNH-4-Pyr	3-Azt
2-2109	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
2-2110	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyrd
2-2111	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
2-2112	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pip
2-2113	3,4-diF-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
2-2114	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
2-2115	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
2-2116	3,4-diF-Ph	2-MeNH-4-Pyr	1-Piz
2-2117	3,4-diF-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
2-2118	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pyr
2-2119	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyr
2-2120	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pym
2-2121	3,4-diF-Ph	2-MeNH-4-Pyr	5-Pym
2-2122	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
2-2123	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂
2-2124	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
2-2125	3,4-diF-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
2-2126	3,4-diF-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
2-2127	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
2-2128	3,4-diF-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-2129	3,4-diF-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
2-2130	3,4-diF-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
2-2131	3,4-diF-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
2-2132	3-Cl-Ph	2-MeNH-4-Pyr	H ₂ N-(CH ₂) ₃
2-2133	3-Cl-Ph	2-MeNH-4-Pyr	MeNH-(CH ₂) ₃
2-2134	3-Cl-Ph	2-MeNH-4-Pyr	EtNH-(CH ₂) ₃
2-2135	3-Cl-Ph	2-MeNH-4-Pyr	Me ₂ N-(CH ₂) ₃
2-2136	3-Cl-Ph	2-MeNH-4-Pyr	1-Azt-(CH ₂) ₃
2-2137	3-Cl-Ph	2-MeNH-4-Pyr	1-Pyrd-(CH ₂) ₃
2-2138	3-Cl-Ph	2-MeNH-4-Pyr	1-Pip-(CH ₂) ₃
2-2139	3-Cl-Ph	2-MeNH-4-Pyr	1-Mor-(CH ₂) ₃
2-2140	3-Cl-Ph	2-MeNH-4-Pyr	1-Tmor-(CH ₂) ₃
2-2141	3-Cl-Ph	2-MeNH-4-Pyr	1-Piz-(CH ₂) ₃
2-2142	3-Cl-Ph	2-MeNH-4-Pyr	4-Me-1-Piz-(CH ₂) ₃
2-2143	3-Cl-Ph	2-MeNH-4-Pyr	3-Azt
2-2144	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Azt
2-2145	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyrd
2-2146	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd
2-2147	3-Cl-Ph	2-MeNH-4-Pyr	4-Pip
2-2148	3-Cl-Ph	2-MeNH-4-Pyr	4-(3,4-deH-Pip)
2-2149	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-Pip
2-2150	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-(3,4-deH-Pip)
2-2151	3-Cl-Ph	2-MeNH-4-Pyr	1-Piz
2-2152	3-Cl-Ph	2-MeNH-4-Pyr	4-Me-1-Piz
2-2153	3-Cl-Ph	2-MeNH-4-Pyr	4-Pyr
2-2154	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyr
2-2155	3-Cl-Ph	2-MeNH-4-Pyr	4-Pym
2-2156	3-Cl-Ph	2-MeNH-4-Pyr	5-Pym
2-2157	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyrd-CH ₂
2-2158	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-2159	3-Cl-Ph	2-MeNH-4-Pyr	4-Pip-CH ₂
2-2160	3-Cl-Ph	2-MeNH-4-Pyr	1-Me-4-Pip-CH ₂
2-2161	3-Cl-Ph	2-MeNH-4-Pyr	2-Piz-CH ₂
2-2162	3-Cl-Ph	2-MeNH-4-Pyr	4-Pyr-CH ₂
2-2163	3-Cl-Ph	2-MeNH-4-Pyr	3-Pyr-CH ₂
2-2164	3-Cl-Ph	2-MeNH-4-Pyr	4-Pym-CH ₂
2-2165	3-Cl-Ph	2-MeNH-4-Pyr	5-Pym-CH ₂
2-2166	3-Cl-Ph	2-MeNH-4-Pyr	2-Pym-CH ₂
2-2167	Ph	4-Pym	H ₂ N-(CH ₂) ₃
2-2168	Ph	4-Pym	MeNH-(CH ₂) ₃
2-2169	Ph	4-Pym	EtNH-(CH ₂) ₃
2-2170	Ph	4-Pym	Me ₂ N-(CH ₂) ₃
2-2171	Ph	4-Pym	1-Azt-(CH ₂) ₃
2-2172	Ph	4-Pym	1-Pyrd-(CH ₂) ₃
2-2173	Ph	4-Pym	1-Pip-(CH ₂) ₃
2-2174	Ph	4-Pym	1-Mor-(CH ₂) ₃
2-2175	Ph	4-Pym	1-Tmor-(CH ₂) ₃
2-2176	Ph	4-Pym	1-Piz-(CH ₂) ₃
2-2177	Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2178	Ph	4-Pym	3-Azt
2-2179	Ph	4-Pym	1-Me-3-Azt
2-2180	Ph	4-Pym	3-Pyrd
2-2181	Ph	4-Pym	1-Me-3-Pyrd
2-2182	Ph	4-Pym	4-Pip
2-2183	Ph	4-Pym	4-(3,4-deH-Pip)
2-2184	Ph	4-Pym	1-Me-4-Pip
2-2185	Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
2-2186	Ph	4-Pym	1-Piz
2-2187	Ph	4-Pym	4-Me-1-Piz
2-2188	Ph	4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
2-2189	Ph	4-Pym	3-Pyr
2-2190	Ph	4-Pym	4-Pym
2-2191	Ph	4-Pym	5-Pym
2-2192	Ph	4-Pym	3-Pyrd-CH ₂
2-2193	Ph	4-Pym	1-Me-3-Pyrd-CH ₂
2-2194	Ph	4-Pym	4-Pip-CH ₂
2-2195	Ph	4-Pym	1-Me-4-Pip-CH ₂
2-2196	Ph	4-Pym	2-Piz-CH ₂
2-2197	Ph	4-Pym	4-Pyr-CH ₂
2-2198	Ph	4-Pym	3-Pyr-CH ₂
2-2199	Ph	4-Pym	4-Pym-CH ₂
2-2200	Ph	4-Pym	5-Pym-CH ₂
2-2201	Ph	4-Pym	2-Pym-CH ₂
2-2202	4-F-Ph	4-Pym	H ₂ N-(CH ₂) ₃
2-2203	4-F-Ph	4-Pym	MeNH-(CH ₂) ₃
2-2204	4-F-Ph	4-Pym	EtNH-(CH ₂) ₃
2-2205	4-F-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
2-2206	4-F-Ph	4-Pym	1-Azt-(CH ₂) ₃
2-2207	4-F-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
2-2208	4-F-Ph	4-Pym	1-Pip-(CH ₂) ₃
2-2209	4-F-Ph	4-Pym	1-Mor-(CH ₂) ₃
2-2210	4-F-Ph	4-Pym	1-Tmor-(CH ₂) ₃
2-2211	4-F-Ph	4-Pym	1-Piz-(CH ₂) ₃
2-2212	4-F-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2213	4-F-Ph	4-Pym	3-Azt
2-2214	4-F-Ph	4-Pym	1-Me-3-Azt
2-2215	4-F-Ph	4-Pym	3-Pyrd
2-2216	4-F-Ph	4-Pym	1-Me-3-Pyrd
2-2217	4-F-Ph	4-Pym	4-Pip
2-2218	4-F-Ph	4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2219	4-F-Ph	4-Pym	1-Me-4-Pip
2-2220	4-F-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
2-2221	4-F-Ph	4-Pym	1-Piz
2-2222	4-F-Ph	4-Pym	4-Me-1-Piz
2-2223	4-F-Ph	4-Pym	4-Pyr
2-2224	4-F-Ph	4-Pym	3-Pyr
2-2225	4-F-Ph	4-Pym	4-Pym
2-2226	4-F-Ph	4-Pym	5-Pym
2-2227	4-F-Ph	4-Pym	3-Pyrd-CH ₂
2-2228	4-F-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
2-2229	4-F-Ph	4-Pym	4-Pip-CH ₂
2-2230	4-F-Ph	4-Pym	1-Me-4-Pip-CH ₂
2-2231	4-F-Ph	4-Pym	2-Piz-CH ₂
2-2232	4-F-Ph	4-Pym	4-PyT-CH ₂
2-2233	4-F-Ph	4-Pym	3-PyT-CH ₂
2-2234	4-F-Ph	4-Pym	4-Pym-CH ₂
2-2235	4-F-Ph	4-Pym	5-Pym-CH ₂
2-2236	4-F-Ph	4-Pym	2-Pym-CH ₂
2-2237	3-F-Ph	4-Pym	H ₂ N-(CH ₂) ₃
2-2238	3-F-Ph	4-Pym	MeNH-(CH ₂) ₃
2-2239	3-F-Ph	4-Pym	EtNH-(CH ₂) ₃
2-2240	3-F-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
2-2241	3-F-Ph	4-Pym	1-Azt-(CH ₂) ₃
2-2242	3-F-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
2-2243	3-F-Ph	4-Pym	1-Pip-(CH ₂) ₃
2-2244	3-F-Ph	4-Pym	1-Mor-(CH ₂) ₃
2-2245	3-F-Ph	4-Pym	1-Tmor-(CH ₂) ₃
2-2246	3-F-Ph	4-Pym	1-Piz-(CH ₂) ₃
2-2247	3-F-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2248	3-F-Ph	4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
2-2249	3-F-Ph	4-Pym	1-Me-3-Azt
2-2250	3-F-Ph	4-Pym	3-Pyrd
2-2251	3-F-Ph	4-Pym	1-Me-3-Pyrd
2-2252	3-F-Ph	4-Pym	4-Pip
2-2253	3-F-Ph	4-Pym	4-(3,4-deH-Pip)
2-2254	3-F-Ph	4-Pym	1-Me-4-Pip
2-2255	3-F-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
2-2256	3-F-Ph	4-Pym	1-Piz
2-2257	3-F-Ph	4-Pym	4-Me-1-Piz
2-2258	3-F-Ph	4-Pym	4-Pyr
2-2259	3-F-Ph	4-Pym	3-Pyr
2-2260	3-F-Ph	4-Pym	4-Pym
2-2261	3-F-Ph	4-Pym	5-Pym
2-2262	3-F-Ph	4-Pym	3-Pyrd-CH ₂
2-2263	3-F-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
2-2264	3-F-Ph	4-Pym	4-Pip-CH ₂
2-2265	3-F-Ph	4-Pym	1-Me-4-Pip-CH ₂
2-2266	3-F-Ph	4-Pym	2-Piz-CH ₂
2-2267	3-F-Ph	4-Pym	4-Pyr-CH ₂
2-2268	3-F-Ph	4-Pym	3-Pyr-CH ₂
2-2269	3-F-Ph	4-Pym	4-Pym-CH ₂
2-2270	3-F-Ph	4-Pym	5-Pym-CH ₂
2-2271	3-F-Ph	4-Pym	2-Pym-CH ₂
2-2272	3,4-diF-Ph	4-Pym	H ₂ N-(CH ₂) ₃
2-2273	3,4-diF-Ph	4-Pym	MeNH-(CH ₂) ₃
2-2274	3,4-diF-Ph	4-Pym	EtNH-(CH ₂) ₃
2-2275	3,4-diF-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
2-2276	3,4-diF-Ph	4-Pym	1-Azt-(CH ₂) ₃
2-2277	3,4-diF-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
2-2278	3,4-diF-Ph	4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-2279	3,4-diF-Ph	4-Pym	1-Mor-(CH ₂) ₃
2-2280	3,4-diF-Ph	4-Pym	1-Tmor-(CH ₂) ₃
2-2281	3,4-diF-Ph	4-Pym	1-Piz-(CH ₂) ₃
2-2282	3,4-diF-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2283	3,4-diF-Ph	4-Pym	3-Azt
2-2284	3,4-diF-Ph	4-Pym	1-Me-3-Azt
2-2285	3,4-diF-Ph	4-Pym	3-Pyrd
2-2286	3,4-diF-Ph	4-Pym	1-Me-3-Pyrd
2-2287	3,4-diF-Ph	4-Pym	4-Pip
2-2288	3,4-diF-Ph	4-Pym	4-(3,4-deH-Pip)
2-2289	3,4-diF-Ph	4-Pym	1-Me-4-Pip
2-2290	3,4-diF-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
2-2291	3,4-diF-Ph	4-Pym	1-Piz
2-2292	3,4-diF-Ph	4-Pym	4-Me-1-Piz
2-2293	3,4-diF-Ph	4-Pym	4-Pyr
2-2294	3,4-diF-Ph	4-Pym	3-Pyr
2-2295	3,4-diF-Ph	4-Pym	4-Pym
2-2296	3,4-diF-Ph	4-Pym	5-Pym
2-2297	3,4-diF-Ph	4-Pym	3-Pyrd-CH ₂
2-2298	3,4-diF-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
2-2299	3,4-diF-Ph	4-Pym	4-Pip-CH ₂
2-2300	3,4-diF-Ph	4-Pym	1-Me-4-Pip-CH ₂
2-2301	3,4-diF-Ph	4-Pym	2-Piz-CH ₂
2-2302	3,4-diF-Ph	4-Pym	4-Pyr-CH ₂
2-2303	3,4-diF-Ph	4-Pym	3-Pyr-CH ₂
2-2304	3,4-diF-Ph	4-Pym	4-Pym-CH ₂
2-2305	3,4-diF-Ph	4-Pym	5-Pym-CH ₂
2-2306	3,4-diF-Ph	4-Pym	2-Pym-CH ₂
2-2307	3-Cl-Ph	4-Pym	H ₂ N-(CH ₂) ₃
2-2308	3-Cl-Ph	4-Pym	MeNH-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-2309	3-Cl-Ph	4-Pym	EtNH-(CH ₂) ₃
2-2310	3-Cl-Ph	4-Pym	Me ₂ N-(CH ₂) ₃
2-2311	3-Cl-Ph	4-Pym	1-Azt-(CH ₂) ₃
2-2312	3-Cl-Ph	4-Pym	1-Pyrd-(CH ₂) ₃
2-2313	3-Cl-Ph	4-Pym	1-Pip-(CH ₂) ₃
2-2314	3-Cl-Ph	4-Pym	1-Mor-(CH ₂) ₃
2-2315	3-Cl-Ph	4-Pym	1-Tmor-(CH ₂) ₃
2-2316	3-Cl-Ph	4-Pym	1-Piz-(CH ₂) ₃
2-2317	3-Cl-Ph	4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2318	3-Cl-Ph	4-Pym	3-Azt
2-2319	3-Cl-Ph	4-Pym	1-Me-3-Azt
2-2320	3-Cl-Ph	4-Pym	3-Pyrd
2-2321	3-Cl-Ph	4-Pym	1-Me-3-Pyrd
2-2322	3-Cl-Ph	4-Pym	4-Pip
2-2323	3-Cl-Ph	4-Pym	4-(3,4-deH-Pip)
2-2324	3-Cl-Ph	4-Pym	1-Me-4-Pip
2-2325	3-Cl-Ph	4-Pym	1-Me-4-(3,4-deH-Pip)
2-2326	3-Cl-Ph	4-Pym	1-Piz
2-2327	3-Cl-Ph	4-Pym	4-Me-1-Piz
2-2328	3-Cl-Ph	4-Pym	4-Pyr
2-2329	3-Cl-Ph	4-Pym	3-Pyr
2-2330	3-Cl-Ph	4-Pym	4-Pym
2-2331	3-Cl-Ph	4-Pym	5-Pym
2-2332	3-Cl-Ph	4-Pym	3-Pyrd-CH ₂
2-2333	3-Cl-Ph	4-Pym	1-Me-3-Pyrd-CH ₂
2-2334	3-Cl-Ph	4-Pym	4-Pip-CH ₂
2-2335	3-Cl-Ph	4-Pym	1-Me-4-Pip-CH ₂
2-2336	3-Cl-Ph	4-Pym	2-Piz-CH ₂
2-2337	3-Cl-Ph	4-Pym	4-Pyr-CH ₂
2-2338	3-Cl-Ph	4-Pym	3-Pyr-CH ₂

Compound No.	R ¹	R ²	R ³
2-2339	3-Cl-Ph	4-Pym	4-Pym-CH ₂
2-2340	3-Cl-Ph	4-Pym	5-Pym-CH ₂
2-2341	3-Cl-Ph	4-Pym	2-Pym-CH ₂
2-2342	Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
2-2343	Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
2-2344	Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
2-2345	Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
2-2346	Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
2-2347	Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
2-2348	Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
2-2349	Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
2-2350	Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
2-2351	Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
2-2352	Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2353	Ph	2-MeO-4-Pym	3-Azt
2-2354	Ph	2-MeO-4-Pym	1-Me-3-Azt
2-2355	Ph	2-MeO-4-Pym	3-Pyrd
2-2356	Ph	2-MeO-4-Pym	1-Me-3-Pyrd
2-2357	Ph	2-MeO-4-Pym	4-Pip
2-2358	Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
2-2359	Ph	2-MeO-4-Pym	1-Me-4-Pip
2-2360	Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
2-2361	Ph	2-MeO-4-Pym	1-Piz
2-2362	Ph	2-MeO-4-Pym	4-Me-1-Piz
2-2363	Ph	2-MeO-4-Pym	4-Pyr
2-2364	Ph	2-MeO-4-Pym	3-Pyr
2-2365	Ph	2-MeO-4-Pym	4-Pym
2-2366	Ph	2-MeO-4-Pym	5-Pym
2-2367	Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
2-2368	Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂

Compound No.	R ¹	R ²	R ³
2-2369	Ph	2-MeO-4-Pym	4-Pip-CH ₂
2-2370	Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
2-2371	Ph	2-MeO-4-Pym	2-Piz-CH ₂
2-2372	Ph	2-MeO-4-Pym	4-Pyr-CH ₂
2-2373	Ph	2-MeO-4-Pym	3-Pyr-CH ₂
2-2374	Ph	2-MeO-4-Pym	4-Pym-CH ₂
2-2375	Ph	2-MeO-4-Pym	5-Pym-CH ₂
2-2376	Ph	2-MeO-4-Pym	2-Pym-CH ₂
2-2377	4-F-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
2-2378	4-F-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
2-2379	4-F-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
2-2380	4-F-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
2-2381	4-F-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
2-2382	4-F-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
2-2383	4-F-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
2-2384	4-F-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
2-2385	4-F-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
2-2386	4-F-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
2-2387	4-F-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2388	4-F-Ph	2-MeO-4-Pym	3-Azt
2-2389	4-F-Ph	2-MeO-4-Pym	1-Me-3-Azt
2-2390	4-F-Ph	2-MeO-4-Pym	3-Pyrd
2-2391	4-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
2-2392	4-F-Ph	2-MeO-4-Pym	4-Pip
2-2393	4-F-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
2-2394	4-F-Ph	2-MeO-4-Pym	1-Me-4-Pip
2-2395	4-F-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
2-2396	4-F-Ph	2-MeO-4-Pym	1-Piz
2-2397	4-F-Ph	2-MeO-4-Pym	4-Me-1-Piz
2-2398	4-F-Ph	2-MeO-4-Pym	4-Pyr

Compound No.	R ¹	R ²	R ³
2-2399	4-F-Ph	2-MeO-4-Pym	3-Pyr
2-2400	4-F-Ph	2-MeO-4-Pym	4-Pym
2-2401	4-F-Ph	2-MeO-4-Pym	5-Pym
2-2402	4-F-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
2-2403	4-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
2-2404	4-F-Ph	2-MeO-4-Pym	4-Pip-CH ₂
2-2405	4-F-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
2-2406	4-F-Ph	2-MeO-4-Pym	2-Piz-CH ₂
2-2407	4-F-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
2-2408	4-F-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
2-2409	4-F-Ph	2-MeO-4-Pym	4-Pym-CH ₂
2-2410	4-F-Ph	2-MeO-4-Pym	5-Pym-CH ₂
2-2411	4-F-Ph	2-MeO-4-Pym	2-Pym-CH ₂
2-2412	3-F-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
2-2413	3-F-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
2-2414	3-F-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
2-2415	3-F-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
2-2416	3-F-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
2-2417	3-F-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
2-2418	3-F-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
2-2419	3-F-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
2-2420	3-F-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
2-2421	3-F-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
2-2422	3-F-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2423	3-F-Ph	2-MeO-4-Pym	3-Azt
2-2424	3-F-Ph	2-MeO-4-Pym	1-Me-3-Azt
2-2425	3-F-Ph	2-MeO-4-Pym	3-Pyrd
2-2426	3-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
2-2427	3-F-Ph	2-MeO-4-Pym	4-Pip
2-2428	3-F-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2429	3-F-Ph	2-MeO-4-Pym	1-Me-4-Pip
2-2430	3-F-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
2-2431	3-F-Ph	2-MeO-4-Pym	1-Piz
2-2432	3-F-Ph	2-MeO-4-Pym	4-Me-1-Piz
2-2433	3-F-Ph	2-MeO-4-Pym	4-Pyr
2-2434	3-F-Ph	2-MeO-4-Pym	3-Pyr
2-2435	3-F-Ph	2-MeO-4-Pym	4-Pym
2-2436	3-F-Ph	2-MeO-4-Pym	5-Pym
2-2437	3-F-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
2-2438	3-F-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
2-2439	3-F-Ph	2-MeO-4-Pym	4-Pip-CH ₂
2-2440	3-F-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
2-2441	3-F-Ph	2-MeO-4-Pym	2-Piz-CH ₂
2-2442	3-F-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
2-2443	3-F-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
2-2444	3-F-Ph	2-MeO-4-Pym	4-Pym-CH ₂
2-2445	3-F-Ph	2-MeO-4-Pym	5-Pym-CH ₂
2-2446	3-F-Ph	2-MeO-4-Pym	2-Pym-CH ₂
2-2447	3,4-diF-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
2-2448	3,4-diF-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
2-2449	3,4-diF-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
2-2450	3,4-diF-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
2-2451	3,4-diF-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
2-2452	3,4-diF-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
2-2453	3,4-diF-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃
2-2454	3,4-diF-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₃
2-2455	3,4-diF-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
2-2456	3,4-diF-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
2-2457	3,4-diF-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2458	3,4-diF-Ph	2-MeO-4-Pym	3-Azt

Compound No.	R ¹	R ²	R ³
2-2459	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Azt
2-2460	3,4-diF-Ph	2-MeO-4-Pym	3-Pyrd
2-2461	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
2-2462	3,4-diF-Ph	2-MeO-4-Pym	4-Pip
2-2463	3,4-diF-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
2-2464	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-Pip
2-2465	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
2-2466	3,4-diF-Ph	2-MeO-4-Pym	1-Piz
2-2467	3,4-diF-Ph	2-MeO-4-Pym	4-Me-1-Piz
2-2468	3,4-diF-Ph	2-MeO-4-Pym	4-Pyr
2-2469	3,4-diF-Ph	2-MeO-4-Pym	3-Pyr
2-2470	3,4-diF-Ph	2-MeO-4-Pym	4-Pym
2-2471	3,4-diF-Ph	2-MeO-4-Pym	5-Pym
2-2472	3,4-diF-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
2-2473	3,4-diF-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
2-2474	3,4-diF-Ph	2-MeO-4-Pym	4-Pip-CH ₂
2-2475	3,4-diF-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
2-2476	3,4-diF-Ph	2-MeO-4-Pym	2-Piz-CH ₂
2-2477	3,4-diF-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
2-2478	3,4-diF-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
2-2479	3,4-diF-Ph	2-MeO-4-Pym	4-Pym-CH ₂
2-2480	3,4-diF-Ph	2-MeO-4-Pym	5-Pym-CH ₂
2-2481	3,4-diF-Ph	2-MeO-4-Pym	2-Pym-CH ₂
2-2482	3-Cl-Ph	2-MeO-4-Pym	H ₂ N-(CH ₂) ₃
2-2483	3-Cl-Ph	2-MeO-4-Pym	MeNH-(CH ₂) ₃
2-2484	3-Cl-Ph	2-MeO-4-Pym	EtNH-(CH ₂) ₃
2-2485	3-Cl-Ph	2-MeO-4-Pym	Me ₂ N-(CH ₂) ₃
2-2486	3-Cl-Ph	2-MeO-4-Pym	1-Azt-(CH ₂) ₃
2-2487	3-Cl-Ph	2-MeO-4-Pym	1-Pyrd-(CH ₂) ₃
2-2488	3-Cl-Ph	2-MeO-4-Pym	1-Pip-(CH ₂) ₃

Compound No.	R ¹	R ²	R ³
2-2489	3-Cl-Ph	2-MeO-4-Pym	1-Mor-(CH ₂) ₂
2-2490	3-Cl-Ph	2-MeO-4-Pym	1-Tmor-(CH ₂) ₃
2-2491	3-Cl-Ph	2-MeO-4-Pym	1-Piz-(CH ₂) ₃
2-2492	3-Cl-Ph	2-MeO-4-Pym	4-Me-1-Piz-(CH ₂) ₃
2-2493	3-Cl-Ph	2-MeO-4-Pym	3-Azt
2-2494	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Azt
2-2495	3-Cl-Ph	2-MeO-4-Pym	3-Pyrd
2-2496	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Pyrd
2-2497	3-Cl-Ph	2-MeO-4-Pym	4-Pip
2-2498	3-Cl-Ph	2-MeO-4-Pym	4-(3,4-deH-Pip)
2-2499	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-Pip
2-2500	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-(3,4-deH-Pip)
2-2501	3-Cl-Ph	2-MeO-4-Pym	1-Piz
2-2502	3-Cl-Ph	2-MeO-4-Pym	4-Me-1-Piz
2-2503	3-Cl-Ph	2-MeO-4-Pym	4-Pyr
2-2504	3-Cl-Ph	2-MeO-4-Pym	3-Pyr
2-2505	3-Cl-Ph	2-MeO-4-Pym	4-Pym
2-2506	3-Cl-Ph	2-MeO-4-Pym	5-Pym
2-2507	3-Cl-Ph	2-MeO-4-Pym	3-Pyrd-CH ₂
2-2508	3-Cl-Ph	2-MeO-4-Pym	1-Me-3-Pyrd-CH ₂
2-2509	3-Cl-Ph	2-MeO-4-Pym	4-Pip-CH ₂
2-2510	3-Cl-Ph	2-MeO-4-Pym	1-Me-4-Pip-CH ₂
2-2511	3-Cl-Ph	2-MeO-4-Pym	2-Piz-CH ₂
2-2512	3-Cl-Ph	2-MeO-4-Pym	4-Pyr-CH ₂
2-2513	3-Cl-Ph	2-MeO-4-Pym	3-Pyr-CH ₂
2-2514	3-Cl-Ph	2-MeO-4-Pym	4-Pym-CH ₂
2-2515	3-Cl-Ph	2-MeO-4-Pym	5-Pym-CH ₂
2-2516	3-Cl-Ph	2-MeO-4-Pym	2-Pym-CH ₂
2-2517	4-F-Ph	4-Pyr	H ₂ N-CH ₂ CH=CH
2-2518	4-F-Ph	4-Pyr	MeNH-CH ₂ CH=CH

Compound No.	R ¹	R ²	R ³
2-2519	4-F-Ph	4-Pyr	Me ₂ N-CH ₂ CH=CH
2-2520	4-F-Ph	4-Pyr	3-Pip-CH ₂
2-2521	4-F-Ph	4-Pyr	1-Me-3-Pip-CH ₂
2-2522	4-F-Ph	4-Pyr	2-Me-4-Pip
2-2523	4-F-Ph	4-Pyr	2,2,6,6-tetraMe-4-Pip
2-2524	4-F-Ph	4-Pyr	1-Ac-4-Pip
2-2525	4-F-Ph	4-Pyr	1-Ac-4-(3,4-deH-Pip)
2-2526	4-F-Ph	4-Pyr	4-OH-4-Pip
2-2527	4-F-Ph	4-Pyr	4-OH-1-Me-4-Pip
2-2528	4-F-Ph	4-Pyr	AcNH-(CH ₂) ₃
2-2529	4-F-Ph	4-Pyr	4-NH ₂ -cHx
2-2530	4-F-Ph	4-Pyr	4-Pyr-CH(OH)
2-2531	4-F-Ph	4-Pyr	3-Pyr-CH(OH)
2-2532	4-F-Ph	4-Pyr	2-Pyr-CH(OH)
2-2533	4-F-Ph	4-Pyr	CF ₃ CONH-(CH ₂) ₃
2-2534	4-F-Ph	4-Pyr	BzNH-(CH ₂) ₃
2-2535	4-F-Ph	4-Pyr	2,4,6-triF-BzNH-CH ₂
2-2536	4-F-Ph	4-Pyr	MeSO ₂ NH-(CH ₂) ₃
2-2537	4-F-Ph	4-Pyr	1-NO ₂ (CH ₂) ₂ -4-Pip
2-2538	4-F-Ph	4-Pyr	2,3,5,6-tetraF-4-Pyr
2-2539	4-F-Ph	4-Pyr	3-Qun
2-2540	4-F-Ph	4-Pyr	3-(2,3-deH-Qun)
2-2541	4-F-Ph	4-Pyr	3-ABO
2-2542	4-F-Ph	4-Pyr	8-Me-3-ABO
2-2543	4-F-Ph	4-Pyr	3-(2,3-deH-ABO)
2-2544	4-F-Ph	4-Pyr	8-Me-3-(2,3-deH-ABO)
2-2545	4-F-Ph	4-Pyr	3-ABN
2-2546	4-F-Ph	4-Pyr	9-Me-3-ABN
2-2547	4-F-Ph	4-Pyr	3-(2,3-deH-ABN)
2-2548	4-F-Ph	4-Pyr	9-Me-3-(2,3-deH-ABN)

Compound No.	R ¹	R ²	R ³
2-2549	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂ CH=CH
2-2550	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂ CH=CH
2-2551	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂ CH=CH
2-2552	4-F-Ph	2-NH ₂ -4-Pym	3-Pip-CH ₂
2-2553	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-Pip-CH ₂
2-2554	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-Pip
2-2555	4-F-Ph	2-NH ₂ -4-Pym	2,2,6,6-tetraMe-4-Pip
2-2556	4-F-Ph	2-NH ₂ -4-Pym	1-Ac-4-Pip
2-2557	4-F-Ph	2-NH ₂ -4-Pym	1-Ac-4-(3,4-deH-Pip)
2-2558	4-F-Ph	2-NH ₂ -4-Pym	4-OH-4-Pip
2-2559	4-F-Ph	2-NH ₂ -4-Pym	4-OH-1-Me-4-Pip
2-2560	4-F-Ph	2-NH ₂ -4-Pym	AcNH-(CH ₂) ₂
2-2561	4-F-Ph	2-NH ₂ -4-Pym	4-NH ₂ -cHx
2-2562	4-F-Ph	2-NH ₂ -4-Pym	3-Qun
2-2563	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-Qun)
2-2564	4-F-Ph	2-NH ₂ -4-Pym	3-ABO
2-2565	4-F-Ph	2-NH ₂ -4-Pym	8-Me-3-ABO
2-2566	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-ABO)
2-2567	4-F-Ph	2-NH ₂ -4-Pym	8-Me-3-(2,3-deH-ABO)
2-2568	4-F-Ph	2-NH ₂ -4-Pym	3-ABN
2-2569	4-F-Ph	2-NH ₂ -4-Pym	9-Me-3-ABN
2-2570	4-F-Ph	2-NH ₂ -4-Pym	3-(2,3-deH-ABN)
2-2571	4-F-Ph	2-NH ₂ -4-Pym	9-Me-3-(2,3-deH-ABN)
2-2572	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂ CH=CH
2-2573	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂ CH=CH
2-2574	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂ CH=CH
2-2575	4-F-Ph	2-MeNH-4-Pym	3-Pip-CH ₂
2-2576	4-F-Ph	2-MeNH-4-Pym	1-Me-3-Pip-CH ₂
2-2577	4-F-Ph	2-MeNH-4-Pym	2-Me-4-Pip
2-2578	4-F-Ph	2-MeNH-4-Pym	2,2,6,6-tetraMe-4-Pip

Compound No.	R ¹	R ²	R ³
2-2579	4-F-Ph	2-MeNH-4-Pym	1-Ac-4-Pip
2-2580	4-F-Ph	2-MeNH-4-Pym	1-Ac-4-(3,4-deH-Pip)
2-2581	4-F-Ph	2-MeNH-4-Pym	4-OH-4-Pip
2-2582	4-F-Ph	2-MeNH-4-Pym	4-OH-1-Me-4-Pip
2-2583	4-F-Ph	2-MeNH-4-Pym	AcNH-(CH ₂) ₂
2-2584	4-F-Ph	2-MeNH-4-Pym	4-NH ₂ -cHx
2-2585	4-F-Ph	2-MeNH-4-Pym	3-Qun
2-2586	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-Qun)
2-2587	4-F-Ph	2-MeNH-4-Pym	3-ABO
2-2588	4-F-Ph	2-MeNH-4-Pym	8-Me-3-ABO
2-2589	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-ABO)
2-2590	4-F-Ph	2-MeNH-4-Pym	8-Me-3-(2,3-deH-ABO)
2-2591	4-F-Ph	2-MeNH-4-Pym	3-ABN
2-2592	4-F-Ph	2-MeNH-4-Pym	9-Me-3-ABN
2-2593	4-F-Ph	2-MeNH-4-Pym	3-(2,3-deH-ABN)
2-2594	4-F-Ph	2-MeNH-4-Pym	9-Me-3-(2,3-deH-ABN)
2-2595	4-F-Ph	2-BnNH-4-Pyr	4-Pip
2-2596	4-F-Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
2-2597	4-F-Ph	2-BnNH-4-Pym	4-Pip
2-2598	4-F-Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)
2-2599	4-F-Ph	2-(-Me-BnNH)-4-Pyr	4-Pip
2-2600	4-F-Ph	2-(-Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)
2-2601	4-F-Ph	2-(-Me-BnNH)-4-Pym	4-Pip
2-2602	4-F-Ph	2-(-Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
2-2603	3-Cl-Ph	2-BnNH-4-Pyr	4-Pip
2-2604	3-Cl-Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
2-2605	3-Cl-Ph	2-BnNH-4-Pym	4-Pip
2-2606	3-Cl-Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2607	3-Cl-Ph	2-(-Me-BnNH)-4-Pyr	4-Pip
2-2608	3-Cl-Ph	2-(-Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)
2-2609	3-Cl-Ph	2-(-Me-BnNH)-4-Pym	4-Pip
2-2610	3-Cl-Ph	2-(-Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
2-2611	3-CF ₃ -Ph	2-BnNH-4-Pyr	4-Pip
2-2612	3-CF ₃ -Ph	2-BnNH-4-Pyr	4-(3,4-deH-Pip)
2-2613	3-CF ₃ -Ph	2-BnNH-4-Pym	4-Pip
2-2614	3-CF ₃ -Ph	2-BnNH-4-Pym	4-(3,4-deH-Pip)
2-2615	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pyr	4-Pip
2-2616	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pyr	4-(3,4-deH-Pip)
2-2617	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pym	4-Pip
2-2618	3-CF ₃ -Ph	2-(-Me-BnNH)-4-Pym	4-(3,4-deH-Pip)
2-2619	4-F-Ph	4-Pyr	2-NH ₂ -4-Pym
2-2620	4-F-Ph	4-Pyr	2-MeNH-4-pym
2-2621	4-F-Ph	4-Pyr	2-NH ₂ -4-Pyr
2-2622	4-F-Ph	4-Pyr	2-MeNH-4-pyr
2-2623	4-F-Ph	4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2624	4-F-Ph	4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
2-2625	4-F-Ph	4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
2-2626	4-F-Ph	4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2627	4-F-Ph	4-Pyr	3-(3,4-deH-Pip)
2-2628	4-F-Ph	4-Pyr	1-Me-3-(3,4-deH-Pip)
2-2629	4-F-Ph	4-Pyr	1-Et-4-(3,4-deH-Pip)
2-2630	4-F-Ph	4-Pyr	1-Pr-4-(3,4-deH-Pip)
2-2631	4-F-Ph	4-Pyr	1-Pr-4-Pip
2-2632	4-F-Ph	4-Pyr	1-iPr-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2633	4-F-Ph	4-Pyr	1-iPr-4-Pip
2-2634	4-F-Ph	4-Pyr	1-Bu-4-(3,4-deH-Pip)
2-2635	4-F-Ph	4-Pyr	1-tBu-4-(3,4-deH-Pip)
2-2636	4-F-Ph	4-Pyr	1-Pn-4-(3,4-deH-Pip)
2-2637	4-F-Ph	4-Pyr	1-Hx-4-(3,4-deH-Pip)
2-2638	4-F-Ph	4-Pyr	1-Hp-4-(3,4-deH-Pip)
2-2639	4-F-Ph	4-Pyr	1-Oc-4-(3,4-deH-Pip)
2-2640	4-F-Ph	4-Pyr	1-Nn-4-(3,4-deH-Pip)
2-2641	4-F-Ph	4-Pyr	1-cPr-4-(3,4-deH-Pip)
2-2642	4-F-Ph	4-Pyr	1-cPn-4-(3,4-deH-Pip)
2-2643	4-F-Ph	4-Pyr	1-cHx-4-(3,4-deH-Pip)
2-2644	4-F-Ph	4-Pyr	1-Bn-4-(3,4-deH-Pip)
2-2645	4-F-Ph	4-Pyr	1-Phet-4-(3,4-deH-Pip)
2-2646	4-F-Ph	4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
2-2647	4-F-Ph	4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
2-2648	4-F-Ph	4-Pyr	1-Allyl-4-(3,4-deH-Pip)
2-2649	4-F-Ph	4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
2-2650	4-F-Ph	4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
2-2651	4-F-Ph	4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
2-2652	4-F-Ph	4-Pyr	1,2,2,6,6-pentaMe-4-Pip
2-2653	4-F-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-2654	4-F-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-2655	4-F-Ph	4-Pyr	7-octaH-Ind
2-2656	4-F-Ph	4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
2-2657	4-F-Ph	4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
2-2658	4-F-Ph	4-Pyr	8-octaH-Qui
2-2659	4-F-Ph	4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
2-2660	4-F-Ph	4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
2-2661	4-F-Ph	4-Pyr	2,2-diMe-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2662	4-F-Ph	4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)
2-2663	4-F-Ph	4-Pyr	2,6-diMe-4-(3,4-deH-Pip)
2-2664	4-F-Ph	4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
2-2665	4-F-Ph	4-Pyr	2-Me-4-(3,4-deH-Pip)
2-2666	4-F-Ph	4-Pyr	1,2-diMe-4-(3,4-deH-Pip)
2-2667	4-F-Ph	4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
2-2668	4-F-Ph	4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
2-2669	4-F-Ph	4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
2-2670	4-F-Ph	4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)
2-2671	4-F-Ph	4-Pyr	2-Et-4-(3,4-deH-Pip)
2-2672	4-F-Ph	4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
2-2673	4-F-Ph	4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
2-2674	4-F-Ph	4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
2-2675	4-F-Ph	4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
2-2676	4-F-Ph	4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
2-2677	4-F-Ph	4-Pyr	2-Pr-4-(3,4-deH-Pip)
2-2678	4-F-Ph	4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
2-2679	4-F-Ph	4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
2-2680	4-F-Ph	4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
2-2681	4-F-Ph	4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
2-2682	4-F-Ph	4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
2-2683	4-F-Ph	4-Pyr	2-Bu-4-(3,4-deH-Pip)
2-2684	4-F-Ph	4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
2-2685	4-F-Ph	4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
2-2686	4-F-Ph	4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
2-2687	4-F-Ph	4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
2-2688	4-F-Ph	4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
2-2689	4-F-Ph	4-Pyr	2-Allyl-4-(3,4-deH-Pip)
2-2690	4-F-Ph	4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
2-2691	4-F-Ph	4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2692	4-F-Ph	4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)
2-2693	4-F-Ph	4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
2-2694	4-F-Ph	4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
2-2695	4-F-Ph	4-Pyr	2-Bn-4-(3,4-deH-Pip)
2-2696	4-F-Ph	4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)
2-2697	4-F-Ph	4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
2-2698	4-F-Ph	4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
2-2699	4-F-Ph	4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
2-2700	4-F-Ph	4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)
2-2701	4-F-Ph	4-Pyr	2-Phet-4-(3,4-deH-Pip)
2-2702	4-F-Ph	4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
2-2703	4-F-Ph	4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
2-2704	4-F-Ph	4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
2-2705	4-F-Ph	4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
2-2706	4-F-Ph	4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
2-2707	4-F-Ph	4-Pyr	2-Me-4-(4,5-deH-Pip)
2-2708	4-F-Ph	4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
2-2709	4-F-Ph	4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
2-2710	4-F-Ph	4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
2-2711	4-F-Ph	4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
2-2712	4-F-Ph	4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
2-2713	4-F-Ph	4-Pyr	2-Et-4-(4,5-deH-Pip)
2-2714	4-F-Ph	4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
2-2715	4-F-Ph	4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
2-2716	4-F-Ph	4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
2-2717	4-F-Ph	4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
2-2718	4-F-Ph	4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
2-2719	4-F-Ph	4-Pyr	2-Pr-4-(4,5-deH-Pip)
2-2720	4-F-Ph	4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
2-2721	4-F-Ph	4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2722	4-F-Ph	4-Pyr	1,2-diPr-4-(4,5-deH-Pip)
2-2723	4-F-Ph	4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
2-2724	4-F-Ph	4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
2-2725	4-F-Ph	4-Pyr	2-Bu-4-(4,5-deH-Pip)
2-2726	4-F-Ph	4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)
2-2727	4-F-Ph	4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
2-2728	4-F-Ph	4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
2-2729	4-F-Ph	4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
2-2730	4-F-Ph	4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)
2-2731	4-F-Ph	4-Pyr	2-Allyl-4-(4,5-deH-Pip)
2-2732	4-F-Ph	4-Pyr	1-Me-2-Allyl-4-(4,5-deH-Pip)
2-2733	4-F-Ph	4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
2-2734	4-F-Ph	4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
2-2735	4-F-Ph	4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
2-2736	4-F-Ph	4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
2-2737	4-F-Ph	4-Pyr	2-Bn-4-(4,5-deH-Pip)
2-2738	4-F-Ph	4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
2-2739	4-F-Ph	4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
2-2740	4-F-Ph	4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
2-2741	4-F-Ph	4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
2-2742	4-F-Ph	4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
2-2743	4-F-Ph	4-Pyr	2-Phet-4-(4,5-deH-Pip)
2-2744	4-F-Ph	4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
2-2745	4-F-Ph	4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
2-2746	4-F-Ph	4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
2-2747	4-F-Ph	4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
2-2748	4-F-Ph	4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
2-2749	4-F-Ph	2-NH ₂ -4-Pym	H ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2750	4-F-Ph	2-NH ₂ -4-Pym	MeNH-CH ₂ C(Me) ₂ CH ₂
2-2751	4-F-Ph	2-NH ₂ -4-Pym	EtNH-CH ₂ C(Me) ₂ CH ₂

Compound No.	R ¹	R ²	R ³
2-2752	4-F-Ph	2-NH ₂ -4-Pym	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2753	4-F-Ph	2-NH ₂ -4-Pym	3-(3,4-deH-Pip)
2-2754	4-F-Ph	2-NH ₂ -4-Pym	1-Me-3-(3,4-deH-Pip)
2-2755	4-F-Ph	2-NH ₂ -4-Pym	1-Et-4-(3,4-deH-Pip)
2-2756	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-4-(3,4-deH-Pip)
2-2757	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-4-Pip
2-2758	4-F-Ph	2-NH ₂ -4-Pym	1-iPr-4-(3,4-deH-Pip)
2-2759	4-F-Ph	2-NH ₂ -4-Pym	1-iPr-4-Pip
2-2760	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-4-(3,4-deH-Pip)
2-2761	4-F-Ph	2-NH ₂ -4-Pym	1-tBu-4-(3,4-deH-Pip)
2-2762	4-F-Ph	2-NH ₂ -4-Pym	1-Pn-4-(3,4-deH-Pip)
2-2763	4-F-Ph	2-NH ₂ -4-Pym	1-Hx-4-(3,4-deH-Pip)
2-2764	4-F-Ph	2-NH ₂ -4-Pym	1-Hp-4-(3,4-deH-Pip)
2-2765	4-F-Ph	2-NH ₂ -4-Pym	1-Oc-4-(3,4-deH-Pip)
2-2766	4-F-Ph	2-NH ₂ -4-Pym	1-Nn-4-(3,4-deH-Pip)
2-2767	4-F-Ph	2-NH ₂ -4-Pym	1-cPr-4-(3,4-deH-Pip)
2-2768	4-F-Ph	2-NH ₂ -4-Pym	1-cPn-4-(3,4-deH-Pip)
2-2769	4-F-Ph	2-NH ₂ -4-Pym	1-cHx-4-(3,4-deH-Pip)
2-2770	4-F-Ph	2-NH ₂ -4-Pym	1-Bn-4-(3,4-deH-Pip)
2-2771	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-4-(3,4-deH-Pip)
2-2772	4-F-Ph	2-NH ₂ -4-Pym	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
2-2773	4-F-Ph	2-NH ₂ -4-Pym	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
2-2774	4-F-Ph	2-NH ₂ -4-Pym	1-Allyl-4-(3,4-deH-Pip)
2-2775	4-F-Ph	2-NH ₂ -4-Pym	1-Propargyl-4-(3,4-deH-Pip)
2-2776	4-F-Ph	2-NH ₂ -4-Pym	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
2-2777	4-F-Ph	2-NH ₂ -4-Pym	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
2-2778	4-F-Ph	2-NH ₂ -4-Pym	1,2,2,6,6-pentaMe-4-Pip
2-2779	4-F-Ph	2-NH ₂ -4-Pym	7-(1,2,3,5,6,8a-hexaH-Ind)
2-2780	4-F-Ph	2-NH ₂ -4-Pym	7-(1,2,3,5,8,8a-hexaH-Ind)

Compound No.	R ¹	R ²	R ³
2-2781	4-F-Ph	2-NH ₂ -4-Pym	7-octaH-Ind
2-2782	4-F-Ph	2-NH ₂ -4-Pym	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
2-2783	4-F-Ph	2-NH ₂ -4-Pym	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
2-2784	4-F-Ph	2-NH ₂ -4-Pym	8-octaH-Qui
2-2785	4-F-Ph	2-NH ₂ -4-Pym	2,2-diMe-4-(3,4-deH-Pip)
2-2786	4-F-Ph	2-NH ₂ -4-Pym	1,2,2-triMe-4-(3,4-deH-Pip)
2-2787	4-F-Ph	2-NH ₂ -4-Pym	2,2-diMe-4-(4,5-deH-Pip)
2-2788	4-F-Ph	2-NH ₂ -4-Pym	1,2,2-triMe-4-(4,5-deH-Pip)
2-2789	4-F-Ph	2-NH ₂ -4-Pym	2,6-diMe-4-(3,4-deH-Pip)
2-2790	4-F-Ph	2-NH ₂ -4-Pym	1,2,6-triMe-4-(3,4-deH-Pip)
2-2791	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-(3,4-deH-Pip)
2-2792	4-F-Ph	2-NH ₂ -4-Pym	1,2-diMe-4-(3,4-deH-Pip)
2-2793	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Me-4-(3,4-deH-Pip)
2-2794	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Me-4-(3,4-deH-Pip)
2-2795	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Me-4-(3,4-deH-Pip)
2-2796	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Me-4-(3,4-deH-Pip)
2-2797	4-F-Ph	2-NH ₂ -4-Pym	2-Et-4-(3,4-deH-Pip)
2-2798	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Et-4-(3,4-deH-Pip)
2-2799	4-F-Ph	2-NH ₂ -4-Pym	1,2-diEt-4-(3,4-deH-Pip)
2-2800	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Et-4-(3,4-deH-Pip)
2-2801	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Et-4-(3,4-deH-Pip)
2-2802	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Et-4-(3,4-deH-Pip)
2-2803	4-F-Ph	2-NH ₂ -4-Pym	2-Pr-4-(3,4-deH-Pip)
2-2804	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Pr-4-(3,4-deH-Pip)
2-2805	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Pr-4-(3,4-deH-Pip)
2-2806	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPr-4-(3,4-deH-Pip)
2-2807	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Pr-4-(3,4-deH-Pip)
2-2808	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Pr-4-(3,4-deH-Pip)
2-2809	4-F-Ph	2-NH ₂ -4-Pym	2-Bu-4-(3,4-deH-Pip)
2-2810	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bu-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2811	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bu-4-(3,4-deH-Pip)
2-2812	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bu-4-(3,4-deH-Pip)
2-2813	4-F-Ph	2-NH ₂ -4-Pym	1,2-diBu-4-(3,4-deH-Pip)
2-2814	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bu-4-(3,4-deH-Pip)
2-2815	4-F-Ph	2-NH ₂ -4-Pym	2-Allyl-4-(3,4-deH-Pip)
2-2816	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Allyl-4-(3,4-deH-Pip)
2-2817	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Allyl-4-(3,4-deH-Pip)
2-2818	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Allyl-4-(3,4-deH-Pip)
2-2819	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Allyl-4-(3,4-deH-Pip)
2-2820	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Allyl-4-(3,4-deH-Pip)
2-2821	4-F-Ph	2-NH ₂ -4-Pym	2-Bn-4-(3,4-deH-Pip)
2-2822	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bn-4-(3,4-deH-Pip)
2-2823	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bn-4-(3,4-deH-Pip)
2-2824	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bn-4-(3,4-deH-Pip)
2-2825	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Bn-4-(3,4-deH-Pip)
2-2826	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bn-4-(3,4-deH-Pip)
2-2827	4-F-Ph	2-NH ₂ -4-Pym	2-Phet-4-(3,4-deH-Pip)
2-2828	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Phet-4-(3,4-deH-Pip)
2-2829	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Phet-4-(3,4-deH-Pip)
2-2830	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Phet-4-(3,4-deH-Pip)
2-2831	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Phet-4-(3,4-deH-Pip)
2-2832	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPhet-4-(3,4-deH-Pip)
2-2833	4-F-Ph	2-NH ₂ -4-Pym	2-Me-4-(4,5-deH-Pip)
2-2834	4-F-Ph	2-NH ₂ -4-Pym	1,2-diMe-4-(4,5-deH-Pip)
2-2835	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Me-4-(4,5-deH-Pip)
2-2836	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Me-4-(4,5-deH-Pip)
2-2837	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Me-4-(4,5-deH-Pip)
2-2838	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Me-4-(4,5-deH-Pip)
2-2839	4-F-Ph	2-NH ₂ -4-Pym	2-Et-4-(4,5-deH-Pip)
2-2840	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Et-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2841	4-F-Ph	2-NH ₂ -4-Pym	1,2-diEt-4-(4,5-deH-Pip)
2-2842	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Et-4-(4,5-deH-Pip)
2-2843	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Et-4-(4,5-deH-Pip)
2-2844	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Et-4-(4,5-deH-Pip)
2-2845	4-F-Ph	2-NH ₂ -4-Pym	2-Pr-4-(4,5-deH-Pip)
2-2846	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Pr-4-(4,5-deH-Pip)
2-2847	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Pr-4-(4,5-deH-Pip)
2-2848	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPr-4-(4,5-deH-Pip)
2-2849	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Pr-4-(4,5-deH-Pip)
2-2850	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Pr-4-(4,5-deH-Pip)
2-2851	4-F-Ph	2-NH ₂ -4-Pym	2-Bu-4-(4,5-deH-Pip)
2-2852	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bu-4-(4,5-deH-Pip)
2-2853	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bu-4-(4,5-deH-Pip)
2-2854	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bu-4-(4,5-deH-Pip)
2-2855	4-F-Ph	2-NH ₂ -4-Pym	1,2-diBu-4-(4,5-deH-Pip)
2-2856	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bu-4-(4,5-deH-Pip)
2-2857	4-F-Ph	2-NH ₂ -4-Pym	2-Allyl-4-(4,5-deH-Pip)
2-2858	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Allyl-4-(4,5-deH-Pip)
2-2859	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Allyl-4-(4,5-deH-Pip)
2-2860	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Allyl-4-(4,5-deH-Pip)
2-2861	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Allyl-4-(4,5-deH-Pip)
2-2862	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Allyl-4-(4,5-deH-Pip)
2-2863	4-F-Ph	2-NH ₂ -4-Pym	2-Bn-4-(4,5-deH-Pip)
2-2864	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Bn-4-(4,5-deH-Pip)
2-2865	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Bn-4-(4,5-deH-Pip)
2-2866	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Bn-4-(4,5-deH-Pip)
2-2867	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Bn-4-(4,5-deH-Pip)
2-2868	4-F-Ph	2-NH ₂ -4-Pym	1-Phet-2-Bn-4-(4,5-deH-Pip)
2-2869	4-F-Ph	2-NH ₂ -4-Pym	2-Phet-4-(4,5-deH-Pip)
2-2870	4-F-Ph	2-NH ₂ -4-Pym	1-Me-2-Phet-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2871	4-F-Ph	2-NH ₂ -4-Pym	1-Et-2-Phet-4-(4,5-deH-Pip)
2-2872	4-F-Ph	2-NH ₂ -4-Pym	1-Pr-2-Phet-4-(4,5-deH-Pip)
2-2873	4-F-Ph	2-NH ₂ -4-Pym	1-Bu-2-Phet-4-(4,5-deH-Pip)
2-2874	4-F-Ph	2-NH ₂ -4-Pym	1,2-diPhet-4-(4,5-deH-Pip)
2-2875	4-F-Ph	2-MeNH-4-Pym	H ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2876	4-F-Ph	2-MeNH-4-Pym	MeNH-CH ₂ C(Me) ₂ CH ₂
2-2877	4-F-Ph	2-MeNH-4-Pym	EtNH-CH ₂ C(Me) ₂ CH ₂
2-2878	4-F-Ph	2-MeNH-4-Pym	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
2-2879	4-F-Ph	2-MeNH-4-Pym	3-(3,4-deH-Pip)
2-2880	4-F-Ph	2-MeNH-4-Pym	1-Me-3-(3,4-deH-Pip)
2-2881	4-F-Ph	2-MeNH-4-Pym	1-Et-4-(3,4-deH-Pip)
2-2882	4-F-Ph	2-MeNH-4-Pym	1-Pr-4-(3,4-deH-Pip)
2-2883	4-F-Ph	2-MeNH-4-Pym	1-Pr-4-Pip
2-2884	4-F-Ph	2-MeNH-4-Pym	1-iPr-4-(3,4-deH-Pip)
2-2885	4-F-Ph	2-MeNH-4-Pym	1-iPr-4-Pip
2-2886	4-F-Ph	2-MeNH-4-Pym	1-Bu-4-(3,4-deH-Pip)
2-2887	4-F-Ph	2-MeNH-4-Pym	1-tBu-4-(3,4-deH-Pip)
2-2888	4-F-Ph	2-MeNH-4-Pym	1-Pn-4-(3,4-deH-Pip)
2-2889	4-F-Ph	2-MeNH-4-Pym	1-Hx-4-(3,4-deH-Pip)
2-2890	4-F-Ph	2-MeNH-4-Pym	1-Hp-4-(3,4-deH-Pip)
2-2891	4-F-Ph	2-MeNH-4-Pym	1-Oc-4-(3,4-deH-Pip)
2-2892	4-F-Ph	2-MeNH-4-Pym	1-Nn-4-(3,4-deH-Pip)
2-2893	4-F-Ph	2-MeNH-4-Pym	1-cPr-4-(3,4-deH-Pip)
2-2894	4-F-Ph	2-MeNH-4-Pym	1-cPn-4-(3,4-deH-Pip)
2-2895	4-F-Ph	2-MeNH-4-Pym	1-cHx-4-(3,4-deH-Pip)
2-2896	4-F-Ph	2-MeNH-4-Pym	1-Bn-4-(3,4-deH-Pip)
2-2897	4-F-Ph	2-MeNH-4-Pym	1-Phet-4-(3,4-deH-Pip)
2-2898	4-F-Ph	2-MeNH-4-Pym	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
2-2899	4-F-Ph	2-MeNH-4-Pym	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
2-2900	4-F-Ph	2-MeNH-4-Pym	1-Allyl-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2901	4-F-Ph	2-MeNH-4-Pym	1-Propargyl-4-(3,4-deH-Pip)
2-2902	4-F-Ph	2-MeNH-4-Pym	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
2-2903	4-F-Ph	2-MeNH-4-Pym	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
2-2904	4-F-Ph	2-MeNH-4-Pym	1,2,2,6,6-pentaMe-4-Pip
2-2905	4-F-Ph	2-MeNH-4-Pym	7-(1,2,3,5,6,8a-hexaH-Ind)
2-2906	4-F-Ph	2-MeNH-4-Pym	7-(1,2,3,5,8,8a-hexaH-Ind)
2-2907	4-F-Ph	2-MeNH-4-Pym	7-octaH-Ind
2-2908	4-F-Ph	2-MeNH-4-Pym	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
2-2909	4-F-Ph	2-MeNH-4-Pym	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
2-2910	4-F-Ph	2-MeNH-4-Pym	8-octaH-Qui
2-2911	4-F-Ph	2-MeNH-4-Pym	2,2-diMe-4-(3,4-deH-Pip)
2-2912	4-F-Ph	2-MeNH-4-Pym	1,2,2-triMe-4-(3,4-deH-Pip)
2-2913	4-F-Ph	2-MeNH-4-Pym	2,2-diMe-4-(4,5-deH-Pip)
2-2914	4-F-Ph	2-MeNH-4-Pym	1,2,2-triMe-4-(4,5-deH-Pip)
2-2915	4-F-Ph	2-MeNH-4-Pym	2,6-diMe-4-(3,4-deH-Pip)
2-2916	4-F-Ph	2-MeNH-4-Pym	1,2,6-triMe-4-(3,4-deH-Pip)
2-2917	4-F-Ph	2-MeNH-4-Pym	2-Me-4-(3,4-deH-Pip)
2-2918	4-F-Ph	2-MeNH-4-Pym	1,2-diMe-4-(3,4-deH-Pip)
2-2919	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Me-4-(3,4-deH-Pip)
2-2920	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Me-4-(3,4-deH-Pip)
2-2921	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Me-4-(3,4-deH-Pip)
2-2922	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Me-4-(3,4-deH-Pip)
2-2923	4-F-Ph	2-MeNH-4-Pym	2-Et-4-(3,4-deH-Pip)
2-2924	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Et-4-(3,4-deH-Pip)
2-2925	4-F-Ph	2-MeNH-4-Pym	1,2-diEt-4-(3,4-deH-Pip)
2-2926	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Et-4-(3,4-deH-Pip)
2-2927	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Et-4-(3,4-deH-Pip)
2-2928	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Et-4-(3,4-deH-Pip)
2-2929	4-F-Ph	2-MeNH-4-Pym	2-Pr-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2930	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Pr-4-(3,4-deH-Pip)
2-2931	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Pr-4-(3,4-deH-Pip)
2-2932	4-F-Ph	2-MeNH-4-Pym	1,2-diPr-4-(3,4-deH-Pip)
2-2933	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Pr-4-(3,4-deH-Pip)
2-2934	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Pr-4-(3,4-deH-Pip)
2-2935	4-F-Ph	2-MeNH-4-Pym	2-Bu-4-(3,4-deH-Pip)
2-2936	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bu-4-(3,4-deH-Pip)
2-2937	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bu-4-(3,4-deH-Pip)
2-2938	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bu-4-(3,4-deH-Pip)
2-2939	4-F-Ph	2-MeNH-4-Pym	1,2-diBu-4-(3,4-deH-Pip)
2-2940	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bu-4-(3,4-deH-Pip)
2-2941	4-F-Ph	2-MeNH-4-Pym	2-Allyl-4-(3,4-deH-Pip)
2-2942	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Allyl-4-(3,4-deH-Pip)
2-2943	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Allyl-4-(3,4-deH-Pip)
2-2944	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Allyl-4-(3,4-deH-Pip)
2-2945	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Allyl-4-(3,4-deH-Pip)
2-2946	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Allyl-4-(3,4-deH-Pip)
2-2947	4-F-Ph	2-MeNH-4-Pym	2-Bn-4-(3,4-deH-Pip)
2-2948	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bn-4-(3,4-deH-Pip)
2-2949	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bn-4-(3,4-deH-Pip)
2-2950	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bn-4-(3,4-deH-Pip)
2-2951	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Bn-4-(3,4-deH-Pip)
2-2952	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bn-4-(3,4-deH-Pip)
2-2953	4-F-Ph	2-MeNH-4-Pym	2-Phet-4-(3,4-deH-Pip)
2-2954	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Phet-4-(3,4-deH-Pip)
2-2955	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Phet-4-(3,4-deH-Pip)
2-2956	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Phet-4-(3,4-deH-Pip)
2-2957	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Phet-4-(3,4-deH-Pip)
2-2958	4-F-Ph	2-MeNH-4-Pym	1,2-diPhet-4-(3,4-deH-Pip)
2-2959	4-F-Ph	2-MeNH-4-Pym	2-Me-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2960	4-F-Ph	2-MeNH-4-Pym	1,2-diMe-4-(4,5-deH-Pip)
2-2961	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Me-4-(4,5-deH-Pip)
2-2962	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Me-4-(4,5-deH-Pip)
2-2963	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Me-4-(4,5-deH-Pip)
2-2964	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Me-4-(4,5-deH-Pip)
2-2965	4-F-Ph	2-MeNH-4-Pym	2-Et-4-(4,5-deH-Pip)
2-2966	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Et-4-(4,5-deH-Pip)
2-2967	4-F-Ph	2-MeNH-4-Pym	1,2-diEt-4-(4,5-deH-Pip)
2-2968	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Et-4-(4,5-deH-Pip)
2-2969	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Et-4-(4,5-deH-Pip)
2-2970	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Et-4-(4,5-deH-Pip)
2-2971	4-F-Ph	2-MeNH-4-Pym	2-Pr-4-(4,5-deH-Pip)
2-2972	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Pr-4-(4,5-deH-Pip)
2-2973	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Pr-4-(4,5-deH-Pip)
2-2974	4-F-Ph	2-MeNH-4-Pym	1,2-diPr-4-(4,5-deH-Pip)
2-2975	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Pr-4-(4,5-deH-Pip)
2-2976	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Pr-4-(4,5-deH-Pip)
2-2977	4-F-Ph	2-MeNH-4-Pym	2-Bu-4-(4,5-deH-Pip)
2-2978	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bu-4-(4,5-deH-Pip)
2-2979	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bu-4-(4,5-deH-Pip)
2-2980	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bu-4-(4,5-deH-Pip)
2-2981	4-F-Ph	2-MeNH-4-Pym	1,2-diBu-4-(4,5-deH-Pip)
2-2982	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bu-4-(4,5-deH-Pip)
2-2983	4-F-Ph	2-MeNH-4-Pym	2-Allyl-4-(4,5-deH-Pip)
2-2984	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Allyl-4-(4,5-deH-Pip)
2-2985	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Allyl-4-(4,5-deH-Pip)
2-2986	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Allyl-4-(4,5-deH-Pip)
2-2987	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Allyl-4-(4,5-deH-Pip)
2-2988	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Allyl-4-(4,5-deH-Pip)
2-2989	4-F-Ph	2-MeNH-4-Pym	2-Bn-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-2990	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Bn-4-(4,5-deH-Pip)
2-2991	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Bn-4-(4,5-deH-Pip)
2-2992	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Bn-4-(4,5-deH-Pip)
2-2993	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Bn-4-(4,5-deH-Pip)
2-2994	4-F-Ph	2-MeNH-4-Pym	1-Phet-2-Bn-4-(4,5-deH-Pip)
2-2995	4-F-Ph	2-MeNH-4-Pym	2-Phet-4-(4,5-deH-Pip)
2-2996	4-F-Ph	2-MeNH-4-Pym	1-Me-2-Phet-4-(4,5-deH-Pip)
2-2997	4-F-Ph	2-MeNH-4-Pym	1-Et-2-Phet-4-(4,5-deH-Pip)
2-2998	4-F-Ph	2-MeNH-4-Pym	1-Pr-2-Phet-4-(4,5-deH-Pip)
2-2999	4-F-Ph	2-MeNH-4-Pym	1-Bu-2-Phet-4-(4,5-deH-Pip)
2-3000	4-F-Ph	2-MeNH-4-Pym	1,2-diPhet-4-(4,5-deH-Pip)
2-3001	4-F-Ph	2-NH ₂ -4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
2-3002	4-F-Ph	2-NH ₂ -4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
2-3003	4-F-Ph	2-NH ₂ -4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
2-3004	4-F-Ph	2-NH ₂ -4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
2-3005	4-F-Ph	2-NH ₂ -4-Pyr	3-(3,4-deH-Pip)
2-3006	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-3-(3,4-deH-Pip)
2-3007	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-4-(3,4-deH-Pip)
2-3008	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-4-(3,4-deH-Pip)
2-3009	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-4-Pip
2-3010	4-F-Ph	2-NH ₂ -4-Pyr	1-iPr-4-(3,4-deH-Pip)
2-3011	4-F-Ph	2-NH ₂ -4-Pyr	1-iPr-4-Pip
2-3012	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-4-(3,4-deH-Pip)
2-3013	4-F-Ph	2-NH ₂ -4-Pyr	1-tBu-4-(3,4-deH-Pip)
2-3014	4-F-Ph	2-NH ₂ -4-Pyr	1-Pn-4-(3,4-deH-Pip)
2-3015	4-F-Ph	2-NH ₂ -4-Pyr	1-Hx-4-(3,4-deH-Pip)
2-3016	4-F-Ph	2-NH ₂ -4-Pyr	1-Hp-4-(3,4-deH-Pip)
2-3017	4-F-Ph	2-NH ₂ -4-Pyr	1-Oc-4-(3,4-deH-Pip)
2-3018	4-F-Ph	2-NH ₂ -4-Pyr	1-Nn-4-(3,4-deH-Pip)
2-3019	4-F-Ph	2-NH ₂ -4-Pyr	1-cPr-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3020	4-F-Ph	2-NH ₂ -4-Pyr	1-cPn-4-(3,4-deH-Pip)
2-3021	4-F-Ph	2-NH ₂ -4-Pyr	1-cHx-4-(3,4-deH-Pip)
2-3022	4-F-Ph	2-NH ₂ -4-Pyr	1-Bn-4-(3,4-deH-Pip)
2-3023	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-4-(3,4-deH-Pip)
2-3024	4-F-Ph	2-NH ₂ -4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
2-3025	4-F-Ph	2-NH ₂ -4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
2-3026	4-F-Ph	2-NH ₂ -4-Pyr	1-Allyl-4-(3,4-deH-Pip)
2-3027	4-F-Ph	2-NH ₂ -4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
2-3028	4-F-Ph	2-NH ₂ -4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
2-3029	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
2-3030	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2,6,6-pentaMe-4-Pip
2-3031	4-F-Ph	2-NH ₂ -4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3032	4-F-Ph	2-NH ₂ -4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3033	4-F-Ph	2-NH ₂ -4-Pyr	7-octaH-Ind
2-3034	4-F-Ph	2-NH ₂ -4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
2-3035	4-F-Ph	2-NH ₂ -4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
2-3036	4-F-Ph	2-NH ₂ -4-Pyr	8-octaH-Qui
2-3037	4-F-Ph	2-NH ₂ -4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
2-3038	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
2-3039	4-F-Ph	2-NH ₂ -4-Pyr	2,2-diMe-4-(4,5-deH-Pip)
2-3040	4-F-Ph	2-NH ₂ -4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)
2-3041	4-F-Ph	2-NH ₂ -4-Pyr	2,6-diMe-4-(3,4-deH-Pip)
2-3042	4-F-Ph	2-NH ₂ -4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
2-3043	4-F-Ph	2-NH ₂ -4-Pyr	2-Me-4-(3,4-deH-Pip)
2-3044	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diMe-4-(3,4-deH-Pip)
2-3045	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
2-3046	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
2-3047	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
2-3048	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3049	4-F-Ph	2-NH ₂ -4-Pyr	2-Et-4-(3,4-deH-Pip)
2-3050	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
2-3051	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
2-3052	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
2-3053	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
2-3054	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
2-3055	4-F-Ph	2-NH ₂ -4-Pyr	2-Pr-4-(3,4-deH-Pip)
2-3056	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
2-3057	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
2-3058	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
2-3059	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
2-3060	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
2-3061	4-F-Ph	2-NH ₂ -4-Pyr	2-Bu-4-(3,4-deH-Pip)
2-3062	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
2-3063	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
2-3064	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
2-3065	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
2-3066	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
2-3067	4-F-Ph	2-NH ₂ -4-Pyr	2-Allyl-4-(3,4-deH-Pip)
2-3068	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
2-3069	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)
2-3070	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)
2-3071	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
2-3072	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
2-3073	4-F-Ph	2-NH ₂ -4-Pyr	2-Bn-4-(3,4-deH-Pip)
2-3074	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)
2-3075	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
2-3076	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
2-3077	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
2-3078	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3079	4-F-Ph	2-NH ₂ -4-Pyr	2-Phet-4-(3,4-deH-Pip)
2-3080	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
2-3081	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
2-3082	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
2-3083	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
2-3084	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
2-3085	4-F-Ph	2-NH ₂ -4-Pyr	2-Me-4-(4,5-deH-Pip)
2-3086	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
2-3087	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
2-3088	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
2-3089	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
2-3090	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
2-3091	4-F-Ph	2-NH ₂ -4-Pyr	2-Et-4-(4,5-deH-Pip)
2-3092	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
2-3093	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
2-3094	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
2-3095	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
2-3096	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
2-3097	4-F-Ph	2-NH ₂ -4-Pyr	2-Pr-4-(4,5-deH-Pip)
2-3098	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
2-3099	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)
2-3100	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPr-4-(4,5-deH-Pip)
2-3101	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
2-3102	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
2-3103	4-F-Ph	2-NH ₂ -4-Pyr	2-Bu-4-(4,5-deH-Pip)
2-3104	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)
2-3105	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
2-3106	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
2-3107	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
2-3108	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3109	4-F-Ph	2-NH ₂ -4-Pyr	2-Allyl-4-(4,5-deH-Pip)
2-3110	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Allyl-4-(4,5-deH-Pip)
2-3111	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
2-3112	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
2-3113	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
2-3114	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
2-3115	4-F-Ph	2-NH ₂ -4-Pyr	2-Bn-4-(4,5-deH-Pip)
2-3116	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
2-3117	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
2-3118	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
2-3119	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
2-3120	4-F-Ph	2-NH ₂ -4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
2-3121	4-F-Ph	2-NH ₂ -4-Pyr	2-Phet-4-(4,5-deH-Pip)
2-3122	4-F-Ph	2-NH ₂ -4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
2-3123	4-F-Ph	2-NH ₂ -4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
2-3124	4-F-Ph	2-NH ₂ -4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
2-3125	4-F-Ph	2-NH ₂ -4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
2-3126	4-F-Ph	2-NH ₂ -4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
2-3127	4-F-Ph	2-MeNH-4-Pyr	H ₂ N-CH ₂ C(Me) ₂ CH ₂
2-3128	4-F-Ph	2-MeNH-4-Pyr	MeNH-CH ₂ C(Me) ₂ CH ₂
2-3129	4-F-Ph	2-MeNH-4-Pyr	EtNH-CH ₂ C(Me) ₂ CH ₂
2-3130	4-F-Ph	2-MeNH-4-Pyr	Me ₂ N-CH ₂ C(Me) ₂ CH ₂
2-3131	4-F-Ph	2-MeNH-4-Pyr	3-(3,4-deH-Pip)
2-3132	4-F-Ph	2-MeNH-4-Pyr	1-Me-3-(3,4-deH-Pip)
2-3133	4-F-Ph	2-MeNH-4-Pyr	1-Et-4-(3,4-deH-Pip)
2-3134	4-F-Ph	2-MeNH-4-Pyr	1-Pr-4-(3,4-deH-Pip)
2-3135	4-F-Ph	2-MeNH-4-Pyr	1-Pr-4-Pip
2-3136	4-F-Ph	2-MeNH-4-Pyr	1-iPr-4-(3,4-deH-Pip)
2-3137	4-F-Ph	2-MeNH-4-Pyr	1-iPr-4-Pip
2-3138	4-F-Ph	2-MeNH-4-Pyr	1-Bu-4-(3,4-deH-Pip)

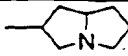
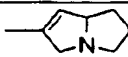
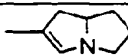
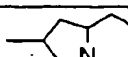
Compound No.	R ¹	R ²	R ³
2-3139	4-F-Ph	2-MeNH-4-Pyr	1-tBu-4-(3,4-deH-Pip)
2-3140	4-F-Ph	2-MeNH-4-Pyr	1-Pn-4-(3,4-deH-Pip)
2-3141	4-F-Ph	2-MeNH-4-Pyr	1-Hx-4-(3,4-deH-Pip)
2-3142	4-F-Ph	2-MeNH-4-Pyr	1-Hp-4-(3,4-deH-Pip)
2-3143	4-F-Ph	2-MeNH-4-Pyr	1-Occ-4-(3,4-deH-Pip)
2-3144	4-F-Ph	2-MeNH-4-Pyr	1-Nn-4-(3,4-deH-Pip)
2-3145	4-F-Ph	2-MeNH-4-Pyr	1-cPr-4-(3,4-deH-Pip)
2-3146	4-F-Ph	2-MeNH-4-Pyr	1-cPn-4-(3,4-deH-Pip)
2-3147	4-F-Ph	2-MeNH-4-Pyr	1-cHx-4-(3,4-deH-Pip)
2-3148	4-F-Ph	2-MeNH-4-Pyr	1-Bn-4-(3,4-deH-Pip)
2-3149	4-F-Ph	2-MeNH-4-Pyr	1-Phet-4-(3,4-deH-Pip)
2-3150	4-F-Ph	2-MeNH-4-Pyr	1-(3-Ph-Pr)-4-(3,4-deH-Pip)
2-3151	4-F-Ph	2-MeNH-4-Pyr	1-(4-Ph-Bu)-4-(3,4-deH-Pip)
2-3152	4-F-Ph	2-MeNH-4-Pyr	1-Allyl-4-(3,4-deH-Pip)
2-3153	4-F-Ph	2-MeNH-4-Pyr	1-Propargyl-4-(3,4-deH-Pip)
2-3154	4-F-Ph	2-MeNH-4-Pyr	2,2,6,6-tetraMe-4-(3,4-deH-Pip)
2-3155	4-F-Ph	2-MeNH-4-Pyr	1,2,2,6,6-pentaMe-4-(3,4-deH-Pip)
2-3156	4-F-Ph	2-MeNH-4-Pyr	1,2,2,6,6-pentaMe-4-Pip
2-3157	4-F-Ph	2-MeNH-4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3158	4-F-Ph	2-MeNH-4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3159	4-F-Ph	2-MeNH-4-Pyr	7-octaH-Ind
2-3160	4-F-Ph	2-MeNH-4-Pyr	8-(1,3,4,6,7,9a-hexaH-2H-Qui)
2-3161	4-F-Ph	2-MeNH-4-Pyr	8-(1,3,4,6,9,9a-hexaH-2H-Qui)
2-3162	4-F-Ph	2-MeNH-4-Pyr	8-octaH-Qui
2-3163	4-F-Ph	2-MeNH-4-Pyr	2,2-diMe-4-(3,4-deH-Pip)
2-3164	4-F-Ph	2-MeNH-4-Pyr	1,2,2-triMe-4-(3,4-deH-Pip)
2-3165	4-F-Ph	2-MeNH-4-Pyr	2,2-diMe-4-(4,5-deH-Pip)
2-3166	4-F-Ph	2-MeNH-4-Pyr	1,2,2-triMe-4-(4,5-deH-Pip)
2-3167	4-F-Ph	2-MeNH-4-Pyr	2,6-diMe-4-(3,4-deH-Pip)

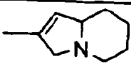
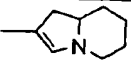
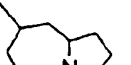

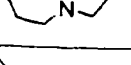
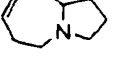
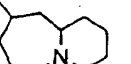
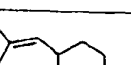
Compound No.	R ¹	R ²	R ³
2-3168	4-F-Ph	2-MeNH-4-Pyr	1,2,6-triMe-4-(3,4-deH-Pip)
2-3169	4-F-Ph	2-MeNH-4-Pyr	2-Me-4-(3,4-deH-Pip)
2-3170	4-F-Ph	2-MeNH-4-Pyr	1,2-diMe-4-(3,4-deH-Pip)
2-3171	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Me-4-(3,4-deH-Pip)
2-3172	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Me-4-(3,4-deH-Pip)
2-3173	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Me-4-(3,4-deH-Pip)
2-3174	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Me-4-(3,4-deH-Pip)
2-3175	4-F-Ph	2-MeNH-4-Pyr	2-Et-4-(3,4-deH-Pip)
2-3176	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Et-4-(3,4-deH-Pip)
2-3177	4-F-Ph	2-MeNH-4-Pyr	1,2-diEt-4-(3,4-deH-Pip)
2-3178	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Et-4-(3,4-deH-Pip)
2-3179	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Et-4-(3,4-deH-Pip)
2-3180	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Et-4-(3,4-deH-Pip)
2-3181	4-F-Ph	2-MeNH-4-Pyr	2-Pr-4-(3,4-deH-Pip)
2-3182	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Pr-4-(3,4-deH-Pip)
2-3183	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Pr-4-(3,4-deH-Pip)
2-3184	4-F-Ph	2-MeNH-4-Pyr	1,2-diPr-4-(3,4-deH-Pip)
2-3185	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Pr-4-(3,4-deH-Pip)
2-3186	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Pr-4-(3,4-deH-Pip)
2-3187	4-F-Ph	2-MeNH-4-Pyr	2-Bu-4-(3,4-deH-Pip)
2-3188	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bu-4-(3,4-deH-Pip)
2-3189	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bu-4-(3,4-deH-Pip)
2-3190	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bu-4-(3,4-deH-Pip)
2-3191	4-F-Ph	2-MeNH-4-Pyr	1,2-diBu-4-(3,4-deH-Pip)
2-3192	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bu-4-(3,4-deH-Pip)
2-3193	4-F-Ph	2-MeNH-4-Pyr	2-Allyl-4-(3,4-deH-Pip)
2-3194	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Allyl-4-(3,4-deH-Pip)
2-3195	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Allyl-4-(3,4-deH-Pip)
2-3196	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Allyl-4-(3,4-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3197	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Allyl-4-(3,4-deH-Pip)
2-3198	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Allyl-4-(3,4-deH-Pip)
2-3199	4-F-Ph	2-MeNH-4-Pyr	2-Bn-4-(3,4-deH-Pip)
2-3200	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bn-4-(3,4-deH-Pip)
2-3201	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bn-4-(3,4-deH-Pip)
2-3202	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bn-4-(3,4-deH-Pip)
2-3203	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Bn-4-(3,4-deH-Pip)
2-3204	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bn-4-(3,4-deH-Pip)
2-3205	4-F-Ph	2-MeNH-4-Pyr	2-Phet-4-(3,4-deH-Pip)
2-3206	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Phet-4-(3,4-deH-Pip)
2-3207	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Phet-4-(3,4-deH-Pip)
2-3208	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Phet-4-(3,4-deH-Pip)
2-3209	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Phet-4-(3,4-deH-Pip)
2-3210	4-F-Ph	2-MeNH-4-Pyr	1,2-diPhet-4-(3,4-deH-Pip)
2-3211	4-F-Ph	2-MeNH-4-Pyr	2-Me-4-(4,5-deH-Pip)
2-3212	4-F-Ph	2-MeNH-4-Pyr	1,2-diMe-4-(4,5-deH-Pip)
2-3213	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Me-4-(4,5-deH-Pip)
2-3214	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Me-4-(4,5-deH-Pip)
2-3215	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Me-4-(4,5-deH-Pip)
2-3216	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Me-4-(4,5-deH-Pip)
2-3217	4-F-Ph	2-MeNH-4-Pyr	2-Et-4-(4,5-deH-Pip)
2-3218	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Et-4-(4,5-deH-Pip)
2-3219	4-F-Ph	2-MeNH-4-Pyr	1,2-diEt-4-(4,5-deH-Pip)
2-3220	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Et-4-(4,5-deH-Pip)
2-3221	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Et-4-(4,5-deH-Pip)
2-3222	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Et-4-(4,5-deH-Pip)
2-3223	4-F-Ph	2-MeNH-4-Pyr	2-Pr-4-(4,5-deH-Pip)
2-3224	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Pr-4-(4,5-deH-Pip)
2-3225	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Pr-4-(4,5-deH-Pip)
2-3226	4-F-Ph	2-MeNH-4-Pyr	1,2-diPr-4-(4,5-deH-Pip)

Compound No.	R ¹	R ²	R ³
2-3227	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Pr-4-(4,5-deH-Pip)
2-3228	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Pr-4-(4,5-deH-Pip)
2-3229	4-F-Ph	2-MeNH-4-Pyr	2-Bu-4-(4,5-deH-Pip)
2-3230	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bu-4-(4,5-deH-Pip)
2-3231	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bu-4-(4,5-deH-Pip)
2-3232	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bu-4-(4,5-deH-Pip)
2-3233	4-F-Ph	2-MeNH-4-Pyr	1,2-diBu-4-(4,5-deH-Pip)
2-3234	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bu-4-(4,5-deH-Pip)
2-3235	4-F-Ph	2-MeNH-4-Pyr	2-Allyl-4-(4,5-deH-Pip)
2-3236	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Allyl-4-(4,5-deH-Pip)
2-3237	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Allyl-4-(4,5-deH-Pip)
2-3238	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Allyl-4-(4,5-deH-Pip)
2-3239	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Allyl-4-(4,5-deH-Pip)
2-3240	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Allyl-4-(4,5-deH-Pip)
2-3241	4-F-Ph	2-MeNH-4-Pyr	2-Bn-4-(4,5-deH-Pip)
2-3242	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Bn-4-(4,5-deH-Pip)
2-3243	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Bn-4-(4,5-deH-Pip)
2-3244	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Bn-4-(4,5-deH-Pip)
2-3245	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Bn-4-(4,5-deH-Pip)
2-3246	4-F-Ph	2-MeNH-4-Pyr	1-Phet-2-Bn-4-(4,5-deH-Pip)
2-3247	4-F-Ph	2-MeNH-4-Pyr	2-Phet-4-(4,5-deH-Pip)
2-3248	4-F-Ph	2-MeNH-4-Pyr	1-Me-2-Phet-4-(4,5-deH-Pip)
2-3249	4-F-Ph	2-MeNH-4-Pyr	1-Et-2-Phet-4-(4,5-deH-Pip)
2-3250	4-F-Ph	2-MeNH-4-Pyr	1-Pr-2-Phet-4-(4,5-deH-Pip)
2-3251	4-F-Ph	2-MeNH-4-Pyr	1-Bu-2-Phet-4-(4,5-deH-Pip)
2-3252	4-F-Ph	2-MeNH-4-Pyr	1,2-diPhet-4-(4,5-deH-Pip)
2-3253	Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3254	Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3255	3-F-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3256	3-F-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)

Compound No.	R ¹	R ²	R ³
2-3257	3-Cl-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3258	3-Cl-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3259	3,4-diF-Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3260	3,4-diF-Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3261	3-CF ₃ -Ph	4-Pyr	7-(1,2,3,5,6,8a-hexaH-Ind)
2-3262	3-CF ₃ -Ph	4-Pyr	7-(1,2,3,5,8,8a-hexaH-Ind)
2-3263	4-F-Ph	4-Pyr	2-Pyrd-CH=CH-
2-3264	4-F-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3265	4-F-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3266	4-F-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3267	4-F-Ph	4-Pyr	2-Pip-CH=CH-
2-3268	4-F-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3269	4-F-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3270	4-F-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
2-3271	Ph	4-Pyr	2-Pyrd-CH=CH-
2-3272	Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3273	Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3274	Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3275	Ph	4-Pyr	2-Pip-CH=CH-
2-3276	Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3277	Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3278	Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
2-3279	3-F-Ph	4-Pyr	2-Pyrd-CH=CH-
2-3280	3-F-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3281	3-F-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3282	3-F-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3283	3-F-Ph	4-Pyr	2-Pip-CH=CH-
2-3284	3-F-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3285	3-F-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3286	3-F-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-

Compound No.	R ¹	R ²	R ³
2-3287	3-Cl-Ph	4-Pyr	2-Pyrd-CH=CH-
2-3288	3-Cl-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3289	3-Cl-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3290	3-Cl-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3291	3-Cl-Ph	4-Pyr	2-Pip-CH=CH-
2-3292	3-Cl-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3293	3-Cl-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3294	3-Cl-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
2-3295	3,4-diF-Ph	4-Pyr	2-Pyrd-CH=CH-
2-3296	3,4-diF-Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3297	3,4-diF-Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3298	3,4-diF-Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3299	3,4-diF-Ph	4-Pyr	2-Pip-CH=CH-
2-3300	3,4-diF-Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3301	3,4-diF-Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3302	3,4-diF-Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
2-3303	3-CF ₃ -Ph	4-Pyr	2-Pyrd-CH=CH-
2-3304	3-CF ₃ -Ph	4-Pyr	1-Me-2-Pyrd-CH=CH-
2-3305	3-CF ₃ -Ph	4-Pyr	1-Et-2-Pyrd-CH=CH-
2-3306	3-CF ₃ -Ph	4-Pyr	1-Pr-2-Pyrd-CH=CH-
2-3307	3-CF ₃ -Ph	4-Pyr	2-Pip-CH=CH-
2-3308	3-CF ₃ -Ph	4-Pyr	1-Me-2-Pip-CH=CH-
2-3309	3-CF ₃ -Ph	4-Pyr	1-Et-2-Pip-CH=CH-
2-3310	3-CF ₃ -Ph	4-Pyr	1-Pr-2-Pip-CH=CH-
2-3311	4-F-Ph	4-Pyr	
2-3312	4-F-Ph	4-Pyr	
2-3313	4-F-Ph	4-Pyr	
2-3314	4-F-Ph	4-Pyr	

Compound No.	R ¹	R ²	R ³
2-3315	4-F-Ph	4-Pyr	
2-3316	4-F-Ph	4-Pyr	
2-3317	4-F-Ph	4-Pyr	
2-3318	4-F-Ph	4-Pyr	
2-3319	4-F-Ph	4-Pyr	
2-3320	4-F-Ph	4-Pyr	
2-3321	4-F-Ph	4-Pyr	
2-3322	4-F-Ph	4-Pyr	

[0093] Of the above compounds in Tables 1 and 2, the following are preferred, that is to say Compounds Nos.: 1-1 to 1-32, 1-35, 1-38, 1-40, 1-42, 1-43, 1-45 to 1-52, 1-65 to 1-67, 1-71 to 1-76, 1-88 to 1-90, 1-94 to 1-126, 1-129, 1-132, 1-134 to 1-137, 1-139 to 1-146, 1-159 to 1-161, 1-165 to 1-170, 1-182 to 1-184, 1-188 to 1-219, 1-222, 1-225, 1-227, 1-229, 1-230, 1-232 to 1-239, 1-252 to 1-254, 1-258 to 1-263, 1-275 to 1-277, 1-281 to 1-312, 1-315, 1-318, 1-320, 1-322, 1-323, 1-325 to 1-332, 1-345 to 1-347, 1-351 to 1-356, 1-368 to 1-370, 1-374 to 1-405, 1-408, 1-411, 1-413, 1-415, 1-416, 1-418 to 1-425, 1-438 to 1-440, 1-444 to 1-449, 1-461 to 1-463, 1-467 to 1-498, 1-501, 1-504, 1-506, 1-508, 1-509, 1-511 to 1-518, 1-531 to 1-533, 1-537 to 1-542, 1-554 to 1-556, 1-560 to 1-591, 1-594, 1-597, 1-599, 1-601, 1-602, 1-604 to 1-611, 1-624 to 1-626, 1-630 to 1-635, 1-647 to 1-649, 1-653 to 1-684, 1-687, 1-690, 1-692, 1-694, 1-695, 1-697 to 1-704, 1-717 to 1-719, 1-723 to 1-728, 1-740 to 1-742, 1-746 to 1-777, 1-780, 1-783, 1-785, 1-787, 1-788, 1-790 to 1-797, 1-810 to 1-812, 1-816 to 1-821, 1-833 to 1-835, 1-839 to 1-870, 1-873, 1-876, 1-878, 1-880, 1-881, 1-883 to 1-890, 1-903 to 1-905, 1-909 to 1-914, 1-926 to 1-928, 1-932 to 1-963, 1-966, 1-969, 1-971, 1-973, 1-974, 1-976 to 1-983, 1-996 to 1-998, 1-1002 to 1-1007, 1-1019 to 1-1021, 1-1025 to 1-1056, 1-1059, 1-1062, 1-1064, 1-1066, 1-1067, 1-1069 to 1-1076, 1-1089 to 1-1091, 1-1095 to 1-1100, 1-1112 to 1-1114, 1-1118 to 1-1149, 1-1152, 1-1155, 1-1157, 1-1159, 1-1160, 1-1162 to 1-1169, 1-1182 to 1-1184, 1-1188 to 1-1193, 1-1205 to 1-1207, 1-1211 to 1-1242, 1-1245, 1-1248, 1-1250, 1-1252, 1-1253, 1-1255 to 1-1262, 1-1275 to 1-1277, 1-1281 to 1-1286, 1-1298 to 1-1300, 1-1304 to 1-1335, 1-1338, 1-1341, 1-1343, 1-1345, 1-1346, 1-1348 to 1-1355, 1-1368 to 1-1370, 1-1374 to 1-1379, 1-1391 to 1-1393, 1-1397 to 1-1405, 1-1407 to 1-1415, 1-1418, 1-1419, 1-1422 to 1-1425, 1-1427, 1-1428, 1-1432 to 1-1440, 1-1442 to 1-1450, 1-1453, 1-1454, 1-1457 to 1-1460, 1-1462, 1-1463, 1-1467 to 1-1475, 1-1477 to 1-1485, 1-1488, 1-1489, 1-1492 to 1-1495, 1-1497, 1-1498, 1-1502 to 1-1510, 1-1512 to 1-1520, 1-1523, 1-1524, 1-1527 to 1-1530, 1-1532, 1-1533, 1-1537 to 1-1545, 1-1547 to 1-1555, 1-1558, 1-1559, 1-1562 to 1-1565, 1-1567, 1-1568, 1-1572 to 1-1580, 1-1582 to 1-1590, 1-1593, 1-1594, 1-1597 to 1-1600, 1-1602, 1-1603, 1-1607 to 1-1615,

1-1617 to 1-1625, 1-1628, 1-1629, 1-1632 to 1-1635, 1-1637, 1-1638, 1-1642 to 1-1650, 1-1652 to 1-1660, 1-1663,
 1-1664, 1-1667 to 1-1670, 1-1672, 1-1673, 1-1677 to 1-1685, 1-1687 to 1-1695, 1-1698, 1-1699, 1-1702 to 1-1705,
 1-1707, 1-1708, 1-1712 to 1-1720, 1-1722 to 1-1730, 1-1733, 1-1734, 1-1737 to 1-1740, 1-1742, 1-1743, 1-1747 to
 1-1755, 1-1757 to 1-1765, 1-1768, 1-1769, 1-1772 to 1-1775, 1-1777, 1-1778, 1-1782 to 1-1790, 1-1792 to 1-1800,
 5 1-1803, 1-1804, 1-1807 to 1-1810, 1-1812, 1-1813, 1-1817 to 1-1825, 1-1827 to 1-1835, 1-1838, 1-1839, 1-1842 to
 1-1845, 1-1847, 1-1848, 1-1852 to 1-1860, 1-1862 to 1-1870, 1-1873, 1-1874, 1-1877 to 1-1880, 1-1882, 1-1883,
 1-1887 to 1-1895, 1-1897 to 1-1905, 1-1908, 1-1909, 1-1912 to 1-1915, 1-1917, 1-1918, 1-1922 to 1-1930, 1-1932 to
 1-1940, 1-1943, 1-1944, 1-1947 to 1-1950, 1-1952, 1-1953, 1-1957 to 1-1965, 1-1967 to 1-1975, 1-1978, 1-1979,
 1-1982 to 1-1985, 1-1987, 1-1988, 1-1992 to 1-2000, 1-2002 to 1-2010, 1-2013, 1-2014, 1-2017 to 1-2020, 1-2022,
 10 1-2023, 1-2027 to 1-2035, 1-2037 to 1-2045, 1-2048, 1-2049, 1-2052 to 1-2055, 1-2057, 1-2058, 1-2062 to 1-2070,
 1-2072 to 1-2080, 1-2083, 1-2084, 1-2087 to 1-2090, 1-2092, 1-2093, 1-2097 to 1-2105, 1-2107 to 1-2115, 1-2118,
 1-2119, 1-2122 to 1-2125, 1-2127, 1-2128, 1-2132 to 1-2140, 1-2142 to 1-2150, 1-2153, 1-2154, 1-2157 to 1-2160,
 1-2162, 1-2163, 1-2167 to 1-2175, 1-2177 to 1-2185, 1-2188, 1-2189, 1-2192 to 1-2195, 1-2197, 1-2198, 1-2202 to
 1-2210, 1-2212 to 1-2220, 1-2223, 1-2224, 1-2227 to 1-2230, 1-2232, 1-2233, 1-2237 to 1-2245, 1-2247 to 1-2255,
 15 1-2258, 1-2259, 1-2262 to 1-2265, 1-2267, 1-2268, 1-2272 to 1-2280, 1-2282 to 1-2290, 1-2293, 1-2294, 1-2297 to
 1-2300, 1-2302, 1-2303, 1-2307 to 1-2315, 1-2317 to 1-2325, 1-2328, 1-2329, 1-2332 to 1-2335, 1-2337, 1-2338,
 1-2342 to 1-2350, 1-2352 to 1-2360, 1-2363, 1-2364, 1-2367 to 1-2370, 1-2372, 1-2373, 1-2377 to 1-2385, 1-2387 to
 1-2395, 1-2398, 1-2399, 1-2402 to 1-2405, 1-2407, 1-2408, 1-2412 to 1-2420, 1-2422 to 1-2430, 1-2433, 1-2434,
 1-2437 to 1-2440, 1-2442, 1-2443, 1-2447 to 1-2455, 1-2457 to 1-2465, 1-2468, 1-2469, 1-2472 to 1-2475, 1-2477,
 20 1-2478, 1-2482 to 1-2490, 1-2492 to 1-2500, 1-2503, 1-2504, 1-2507 to 1-2510, 1-2512, 1-2513, 1-2517 to 1-2529,
 1-2539 to 1-2594, 1-2623 to 1-3322, 2-1 to 2-32, 2-35, 2-38, 2-40, 2-42, 2-43, 2-45 to 2-52, 2-65 to 2-67, 2-71 to 2-76,
 2-88 to 2-90, 2-94 to 2-126, 2-129, 2-132, 2-134 to 2-137, 2-139 to 2-146, 2-159 to 2-161, 2-165 to 2-170, 2-182 to
 2-184, 2-188 to 2-219, 2-222, 2-225, 2-227, 2-229, 2-230, 2-232 to 2-239, 2-252 to 2-254, 2-258 to 2-263, 2-275 to
 2-277, 2-281 to 2-312, 2-315, 2-318, 2-320, 2-322, 2-323, 2-325 to 2-332, 2-345 to 2-347, 2-351 to 2-356, 2-368 to
 25 2-370, 2-374 to 2-405, 2-408, 2-411, 2-413, 2-415, 2-416, 2-418 to 2-425, 2-438 to 2-440, 2-444 to 2-449, 2-461 to
 2-463, 2-467 to 2-498, 2-501, 2-504, 2-506, 2-508, 2-509, 2-511 to 2-518, 2-531 to 2-533, 2-537 to 2-542, 2-554 to
 2-556, 2-560 to 2-591, 2-594, 2-597, 2-599, 2-601, 2-602, 2-604 to 2-611, 2-624 to 2-626, 2-630 to 2-635, 2-647 to
 2-649, 2-653 to 2-684, 2-687, 2-690, 2-692, 2-694, 2-695, 2-697 to 2-704, 2-717 to 2-719, 2-723 to 2-728, 2-740 to
 2-742, 2-746 to 2-777, 2-780, 2-783, 2-785, 2-787, 2-788, 2-790 to 2-797, 2-810 to 2-812, 2-816 to 2-821, 2-833 to
 30 2-835, 2-839 to 2-870, 2-873, 2-876, 2-878, 2-880, 2-881, 2-883 to 2-890, 2-903 to 2-905, 2-909 to 2-914, 2-926 to
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2-2089, 2-2093, 2-2097 to 2-2105, 2-2108 to 2-2115, 2-2118, 2-2119, 2-2122 to 2-2124, 2-2128, 2-2132 to 2-2140, 2-2143 to 2-2150, 2-2153, 2-2154, 2-2157 to 2-2159, 2-2163, 2-2167 to 2-2175, 2-2178 to 2-2185, 2-2188, 2-2189, 2-2192 to 2-2194, 2-2198, 2-2202 to 2-2210, 2-2213 to 2-2220, 2-2223, 2-2224, 2-2227 to 2-2229, 2-2233, 2-2237 to 2-2245, 2-2248 to 2-2255, 2-2258, 2-2259, 2-2262 to 2-2264, 2-2268, 2-2272 to 2-2280, 2-2283 to 2-2290, 2-2293, 2-2294, 2-2297 to 2-2299, 2-2303, 2-2307 to 2-2315, 2-2318 to 2-2325, 2-2328, 2-2329, 2-2332 to 2-2334, 2-2338, 2-2342 to 2-2350, 2-2353 to 2-2360, 2-2363, 2-2364, 2-2367 to 2-2369, 2-2373, 2-2377 to 2-2385, 2-2388 to 2-2395, 2-2398, 2-2399, 2-2402 to 2-2404, 2-2408, 2-2412 to 2-2420, 2-2423 to 2-2430, 2-2433, 2-2434, 2-2437 to 2-2439, 2-2443, 2-2447 to 2-2455, 2-2458 to 2-2465, 2-2468, 2-2469, 2-2472 to 2-2474, 2-2478, 2-2482 to 2-2490, 2-2493 to 2-2500, 2-2503, 2-2504, 2-2507 to 2-2509, 2-2513, 2-2517 to 2-2522, 2-2524 to 2-2526, 2-2528, 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1-1663, 1-1667, 1-1673, 1-1677 to 1-1685, 1-1688, 1-1690, 1-1692 to 1-1695, 1-1698, 1-1702, 1-1706, 1-1712 to 1-1720, 1-1723, 1-1725, 1-1727 to 1-1730, 1-1733, 1-1737, 1-1743, 1-1747 to 1-1755, 1-1758, 1-1760, 1-1762 to 1-1765, 1-1768, 1-1772, 1-1778, 1-1782 to 1-1790, 1-1793, 1-1795, 1-1797 to 1-1800, 1-1803, 1-1807, 1-1813, 1-1817 to 1-1825, 1-1828, 1-1830, 1-1832 to 1-1835, 1-1838, 1-1842, 1-1848, 1-1852 to 1-1860, 1-1863, 1-1865, 1-1867 to 1-1870, 1-1873, 1-1877, 1-1883, 1-1887 to 1-1895, 1-1898, 1-1900, 1-1902 to 1-1905, 1-1908, 1-1912, 1-1918, 1-1922 to 1-1930, 1-1933, 1-1935, 1-1937 to 1-1940, 1-1943, 1-1947, 1-1953, 1-1957 to 1-1965, 1-1968, 1-1970, 1-1972 to 1-1975, 1-1978, 1-1982, 1-1988, 1-1992 to 1-2000, 1-2003, 1-2005, 1-2007 to 1-2010, 1-2013, 1-2017, 1-2023, 1-2027 to 1-2035, 1-2038, 1-2040, 1-2042 to 1-2045, 1-2048, 1-2052, 1-2058, 1-2062 to 1-2070, 1-2073, 1-2075, 1-2077 to 1-2080, 1-2083, 1-2087, 1-2093, 1-2097 to 1-2105, 1-2108, 1-2110, 1-2112 to 1-2115, 1-2118, 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[0094] The most preferred compounds are:

4-(3-aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(3-acetylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(3-methylaminopropyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 [5-(4-fluorophenyl)-4-(pyridin-4-yl)-1*H*-pyrrol-3-yl]-(pyridin-4-yl)methanol,
 2-(1-acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(piperidin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1*H*-pyrrole,
 1-(1-acetyl-piperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-1-[1-(2-nitroethyl)piperidin-4-yl]-2-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-3-yl)-2-(pyridin-4-yl)-1*H*-pyrrole,
 1-(azetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1*H*-pyrrole,
 4-(3-dimethylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-[3-(piperidin-1-yl)propyl]-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-[3-(1-methylpiperazin-4-yl)propyl]-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(2,2,6,6-tetramethylpiperidin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-3-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(quinuclidin-3-yl)-1*H*-pyrrole,
 1-(4-aminocyclohexyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1*H*-pyrrole,
 3-(4-fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1-(piperidin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydro-2,2,6,6-tetramethylpyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(2,2,6,6-tetramethylpiperidin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(1-ethylpiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(1,6-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,2,6,6-pentamethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole,
 4-(3-dimethylamino-1-propen-1-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole.

4-(4-aminobutyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(1-ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole.
 4-(1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole.
 1-(3-dimethylaminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole.
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 4-(6-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(2-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(6-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(2-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2H-quinolizin-8-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,9,9a-hexahydro-2H-quinolizin-8-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1H-pyrrole,
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1H-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1H-pyrrole, and
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1H-pyrrole;

and pharmacologically acceptable salts, esters or other derivatives thereof.

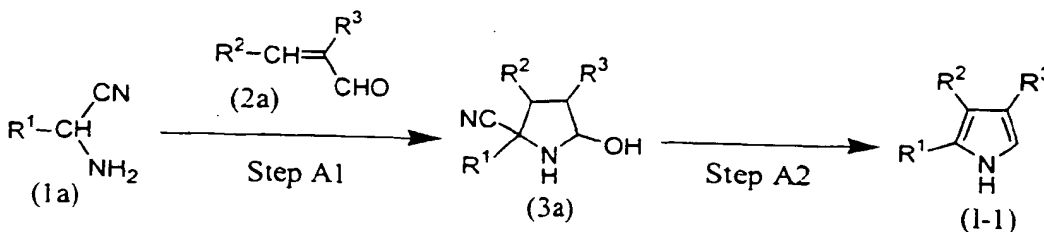
[0095] It should be noted that the specific combination of substituents R¹, R² and R³ shown in the above Tables 1 and 2 can also be applied to the substituents R¹, R² and R³ in formulae (I-2), (I-4) and (I-5) above.

[0096] The compounds of formula (I) and pharmacologically acceptable salts, esters and other derivatives thereof of the present invention can be prepared according to a variety of processes well known in the art for the preparation of compounds of this type, for example as described in the following Methods A to M below.

Method A

[0097] Method A is a method to prepare compounds of formula (I-1) of the present invention.

Reaction Scheme A



[0098] In the above formulae, R¹, R² and R³ are as defined above.

Step A1

[0099] In this Step, a pyrrolidine compound of formula (3a) is prepared by reacting an aminonitrile compound of formula (1a) with an α,β -unsaturated aldehyde compound of formula (2a). This reaction is well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the method

described in EP 0733823, the contents of which are incorporated herein by reference thereto.

[0100] In more detail, this Step is carried out in the presence of a base. There is no particular restriction on the nature of the bases used, and any base commonly used in reactions of this type may equally be used here. Examples of such bases include: alkali metal hydroxides, such as lithium hydroxide, sodium hydroxide and potassium hydroxide; alkali metal hydrides, such as lithium hydride, sodium hydride and potassium hydride; alkali metal amides, such as lithium amide, sodium amide, potassium amide and lithium bis(trimethylsilyl)amide; and alkali metal alkoxides, such as lithium ethoxide, sodium methoxide, sodium ethoxide and potassium t-butoxide. Of these, we prefer the lithium amides.

[0101] The reaction is normally and preferably effected in the presence of a solvent. There is no particular restriction on the nature of the solvent to be employed, provided that it has no adverse effect on the reaction or on the reagents involved and that it can dissolve the reagents, at least to some extent. Examples of suitable solvents include: aliphatic hydrocarbons, such as hexane and heptane; aromatic hydrocarbons, such as benzene, toluene and xylene; ethers, such as diethyl ether, diisopropyl ether, tetrahydrofuran and dioxane; and alcohols, such as methanol, ethanol, propanol, isopropanol and butanol. Of these, we prefer the ethers.

[0102] The reaction can take place over a wide range of temperatures, and the precise reaction temperature is not critical to the invention. The preferred reaction temperature will depend upon such factors as the nature of the solvent, and the starting material or reagent used. However, in general, we find it convenient to carry out the reaction at a temperature of from -78°C to 100°C, more preferably from -78°C to room temperature. The time required for the reaction may also vary widely, depending on many factors, notably the reaction temperature and the nature of the reagents and solvent employed. However, provided that the reaction is effected under the preferred conditions outlined above, a period of from 10 minutes to 30 hours, more preferably from 1 hour to 20 hours, will usually suffice.

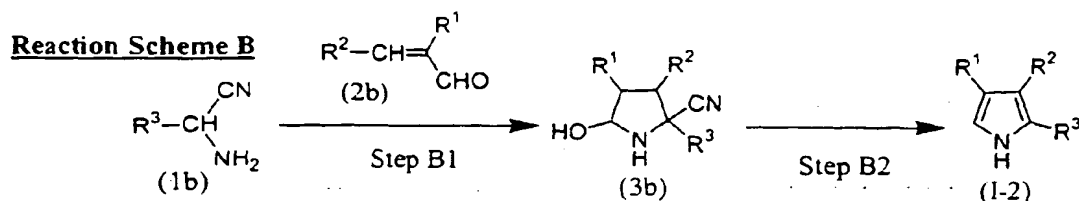
Step A2

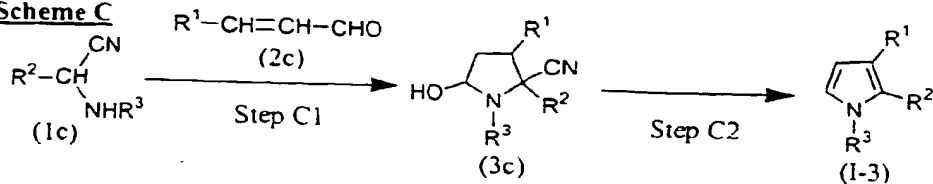
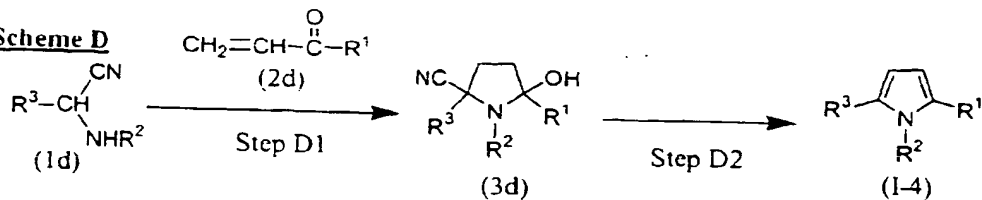
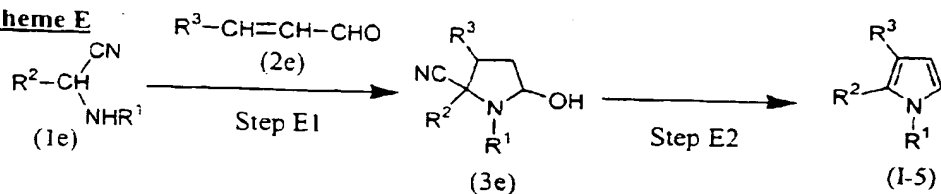
[0103] In this Step, the desired pyrrole derivative of formula (I-1) of the present invention is prepared by the elimination of hydrogen cyanide and water from the compound of formula (3a) prepared in Step A1 above. These reactions are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the methods described in detail in EP 0733823.

[0104] In more detail, this may be achieved by heating the residue obtained by distilling off the solvent from the product of Step A1, or by heating the material obtained by extracting that residue, washing it with water and distilling off the solvent, preferably at a temperature not lower than 100°C, in the presence or absence of a solvent after completion of the reaction of Step A1. The reaction proceeds sufficiently in the absence of a solvent, but, when a solvent is used, the solvent is preferably inert and has a higher boiling point. Examples of suitable solvents include: toluene, xylene, dimethylformamide, dimethylacetamide, dimethyl sulfoxide, diglyme and diphenyl ether.

Method B, Method C, Method D and Method E

[0105] Method B, Method C, Method D and Method E are used to prepare compounds of formulae (I-2), (I-3), (I-4) and (I-5) respectively using the processes outlined in Reaction Schemes B to E below.



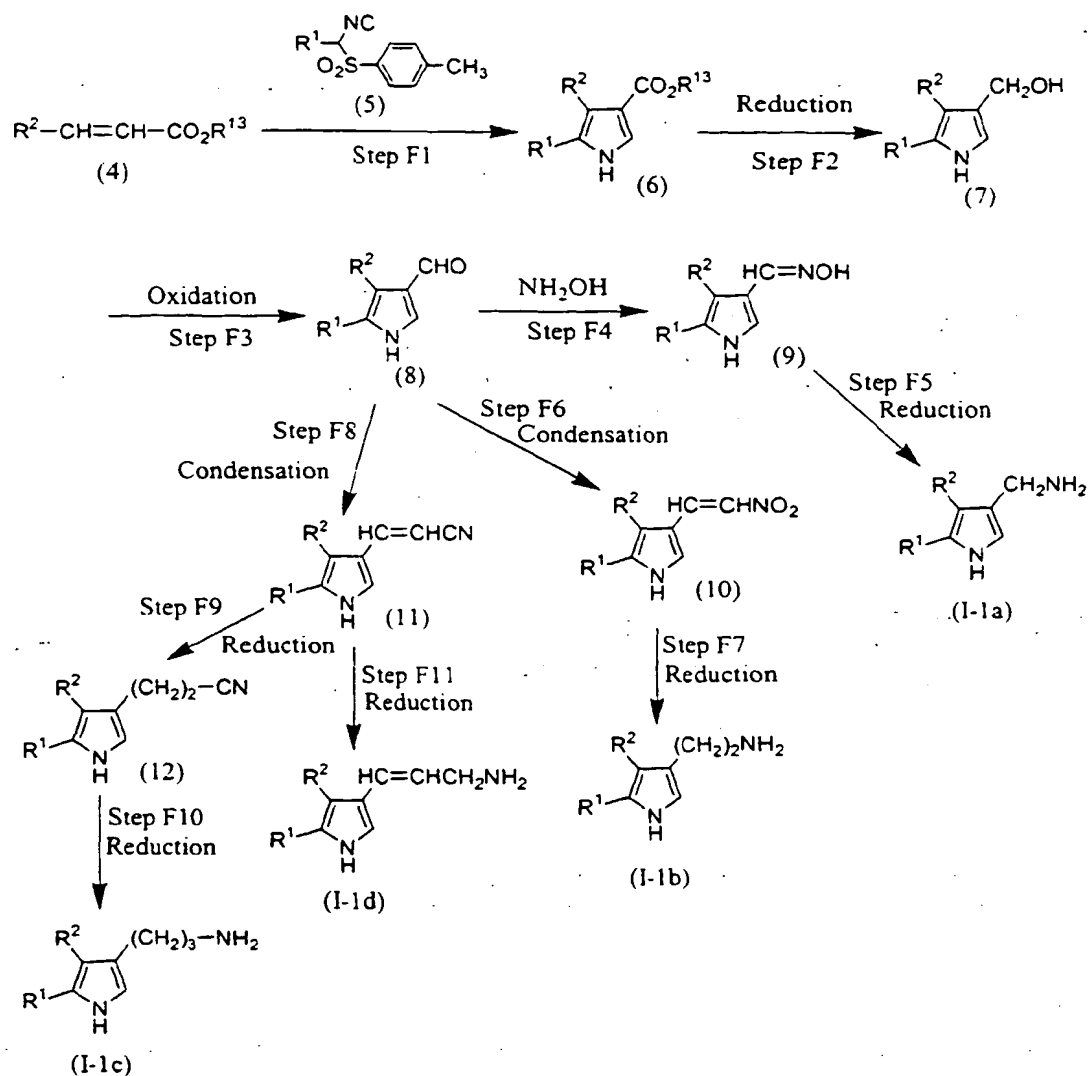
Reaction Scheme C**Reaction Scheme D****Reaction Scheme E**

[0106] In the above formulae, R¹, R² and R³ are as defined above.

[0107] In the above Reaction Schemes Step B1, Step C1, Step D1 and Step E1 are carried out in a similar manner to Step A1 above, and Step B2, Step C2, Step D2 and Step E2 are carried out similarly to Step A2 above.

Method F

[0108] In this method, compounds of formula (I-1) of the present invention wherein R³ is an aminomethyl, aminoethyl, aminopropyl or aminopropenyl group [(I-1a), (I-1b), (I-1c) and (I-1d) respectively] are prepared as shown in Reaction Scheme F.

Reaction Scheme F

[0109] In the above formulae, R^1 and R^2 are as defined above and R^{13} is selected from the group consisting of hydrogen atoms, lower alkyl groups as defined and exemplified above and aralkyl groups as defined and exemplified above.

Step F1

[0110] In this Step, a pyrrole carboxylic acid derivative of formula (6) is prepared by reacting an α,β -unsaturated compound of formula (4) with an isonitrile compound of formula (5). Reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the methods described in detail in R. Di Santo et al., Synthetic Communications, 25(6), pp. 795-802 (1995), the contents of which are incorporated herein by reference thereto.

Step F2

[0111] In this Step, a hydroxymethyl compound of formula (7) is prepared by reducing the pyrrole carboxylic acid derivative of formula (6) prepared in Step F1 above. Reduction reactions of this type are well known in the field of

synthetic organic chemistry and can be carried out using well known techniques, for example according to the method described in detail in R. F. Nystrom et al., J. Am. Chem. Soc., 71, 3245 (1945), the contents of which are incorporated herein by reference thereto, using a reducing agent such as lithium aluminum hydride, lithium borohydride or diisobutyl aluminum hydride.

Step F3

[0112] In this Step, an aldehyde compound of formula (8) is prepared by oxidizing the hydroxymethyl compound of formula (7) prepared in Step F2 above. Oxidation reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the method described in detail in S. Bartel & F. Bohlmann, Tetrahedron Lett., 685 (1985), the contents of which are incorporated herein by reference thereto, using an oxidizing agent such as chromic acid, manganese dioxide or dimethylsulfoxide.

Step F4

[0113] In this Step, an oxime compound of formula (9) is prepared by reacting the aldehyde compound of formula (8) prepared in Step F3 above with hydroxylamine. Dehydration condensation reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques.

Step F5

[0114] In this Step, the desired compound of formula (I-1a) of the present invention is prepared by reducing the oxime compound of formula (9) prepared in Step F4 above. Reduction reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example by catalytic reduction or reduction with diborane, examples of which are disclosed in detail in P. N. Rylander & D. R. Steele, Engelhard Ind. Tech. Bull., 3, 19 (1962), and J. A. Secrist, III & M. W. Logue, J. Org. Chem., 37, 335 (1972), the contents of which are incorporated herein by reference thereto.

Step F6

[0115] In this Step, a nitroolefin compound of formula (10) is prepared by a dehydration condensation (aldol condensation) of the aldehyde compound of formula (8) prepared in Step F3 above and nitromethane. Dehydration condensation reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the method described in detail in D. E. Worall, Org. Synth., I, 413 (1941), the contents of which are incorporated herein by reference thereto.

Step F7

[0116] In this Step, the desired compound of formula (I-1b) of the present invention is prepared by reducing the nitroolefin compound of formula (10) prepared in Step F6 above. Reduction reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the method described in detail in S. I. Murahashi et al., Bull. Chem. Soc. Jpn., 63, 1252 (1990), the contents of which are incorporated herein by reference thereto.

Step F8

[0117] In this Step, a cyanoolefin compound of formula (11) is prepared from the aldehyde compound of formula (8) prepared in Step F3 above. This transformation can be performed, for example, via an aldol condensation reaction similar to that employed in Step F6 above, a Wittig reaction [for example according to the method described in detail in The Organic Chemistry of Phosphorous, Elsevier (1967), the contents of which are incorporated herein by reference thereto], or a Wittig-Horner reaction [for example according to the method described in detail in L. Homer et al., Chem. Ber., 95, 581 (1962), the contents of which are incorporated herein by reference thereto].

Step F9

[0118] In this Step, a cyanoethyl compound of formula (12) is prepared by reduction of the double bond of the cyanoolefin compound of formula (11) prepared in step F8 above. Reduction reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the

method described in detail in S. M. Kerwin et al., J. Org. Chem., 52, 1686 (1987) and T. Hudlicky et al., J. Org. Chem., 52, 4641 (1987), the contents of which are incorporated herein by reference thereto.

Step F10

[0119] In this Step, the desired compound of formula (I-1c) of the present invention is prepared by reducing the cyano group of the cyanoethyl compound of formula (12) prepared in Step F9 above. Reduction reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the methods described in detail in N. M. Yoon & H. C. Brown, J. Am. Chem. Soc., 90, 2927 (1968) and J. Meinwald et al., J. Org. Chem., 29, 2914 (1964), the contents of which are incorporated herein by reference thereto, using lithium aluminum hydride or diisobutyl aluminum hydride.

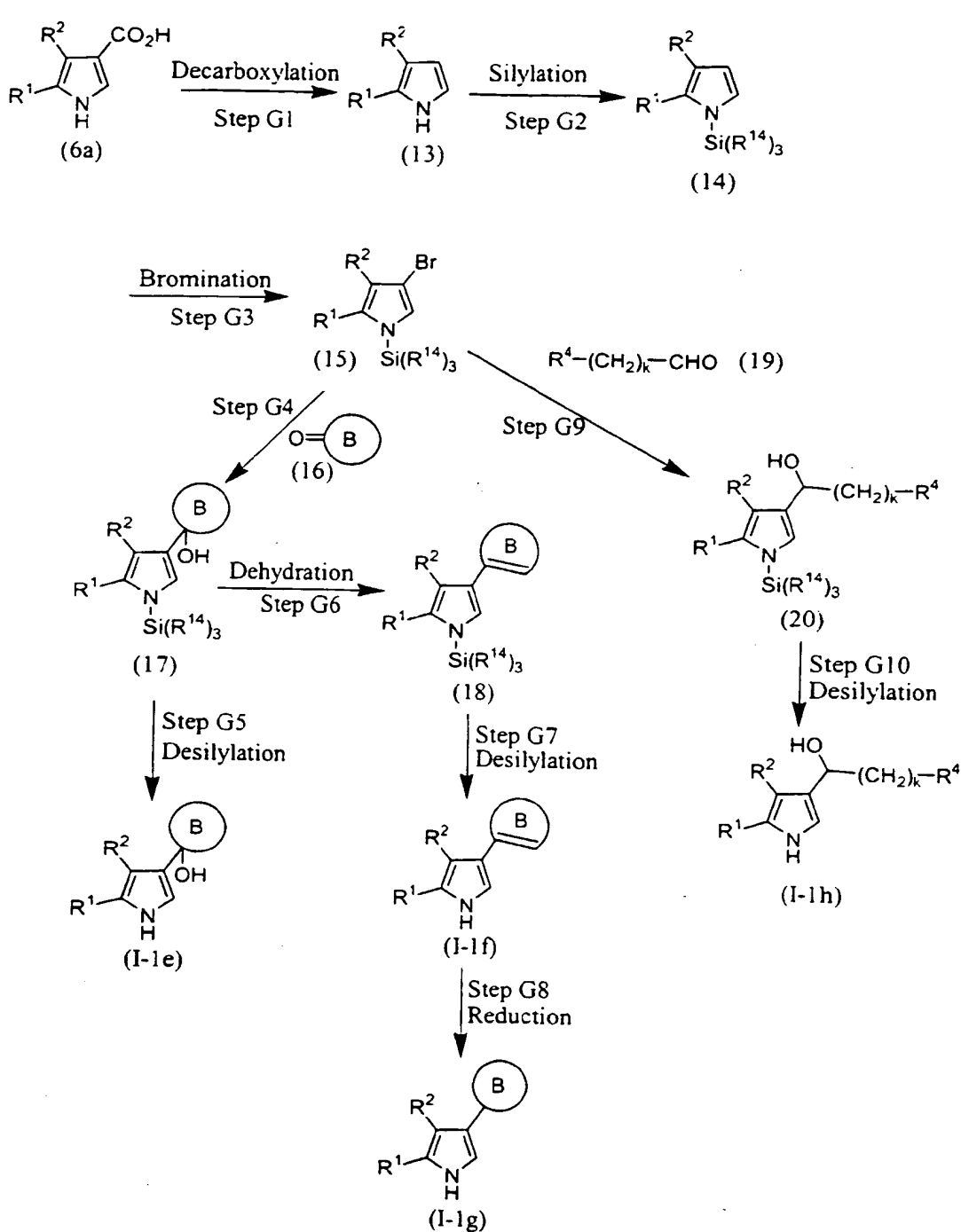
Step F11

[0120] In this Step, the desired compound of formula (I-1d) of the present invention is prepared by reducing the cyano group of the cyanoolefin compound of formula (11) prepared in Step F8 above. This reduction is carried out in a similar manner to that described in Step F10 above.

[0121] Derivatives of the compounds of formula (I-1a), (I-1b), (I-1c) and (I-1d) prepared above in which a nitrogen atom of an amino group present therein is substituted can be synthesized by alkylating, acylating and/or sulfonylating said amino group according to techniques well known in the field of synthetic organic chemistry.

Method G

[0122] In this method, compounds of formula (I-1) of the present invention wherein R³ is a heterocyclyl group which may optionally be substituted with at least one group selected from Substituent group α and Substituent group δ and which is substituted with at least one group selected from Substituent group β and Substituent group γ , as defined above, a heterocyclyl group having at least one nitrogen atom which may be optionally substituted with at least one group selected from Substituent group α and Substituent group δ , as defined above, or a group of the general formula -CH(OH)-(CH₂)_k-R⁴ (wherein R⁴ is as defined above, and the moiety -CH(OH)-(CH₂)_k- is a lower alkylene group as defined above for the group of formula X which is substituted with a hydroxy group, and k is 0 or an integer of from 1 to 5) can be prepared, as shown in Reaction Scheme G below.

Reaction Scheme G

In the above formulae,

R^1 , R^2 , R^4 and k are as defined above,
the substituents R^{14} may be the same or different from one another and each is selected from the group consisting

of hydrogen atoms, lower alkyl groups as defined and exemplified above and aralkyl groups as defined and exemplified above, and

cyclic group B is a heterocycl group corresponding to the heterocycl group which may be optionally substituted with at least one group selected from Substituent group α and Substituent group δ and which is substituted with at least one group selected from Substituent group β and Substituent group γ or the heterocycl group having at least one nitrogen atom, which may be optionally substituted with at least one group selected from Substituent group α and Substituent group δ as defined above in the definition of R^4 .

Step G1

[0123] In this Step, a di-substituted pyrrole compound of formula (13) is prepared by decarboxylating a carboxylic acid compound of formula (6a). Decarboxylation of the compound of formula (6a) can be performed by heating under acidic, alkaline or neutral conditions according to techniques conventionally used in synthetic organic chemistry, for example according to the methods described in detail in N. Yoshida et al., Yakugaku Zasshi, 93(5), 564-598 (1973), the contents of which are incorporated herein by reference thereto.

Steps G2 to G5

[0124] In Step G2, a silyl compound of formula (14) is prepared by silylating the nitrogen atom at the 1-position of the di-substituted pyrrole compound of formula (13) prepared in Step G1 above.

[0125] In Step G3, a bromopyrrole compound of formula (15) is prepared by brominating the silyl compound of formula (14) prepared in Step G2 above with a brominating agent (for example, N-bromosuccinimide, etc.).

[0126] In Step G4, a hydroxyheterocycl compound of formula (17) is prepared by lithiating the compound of formula (15) prepared in Step G3 above and then reacting it with a heterocycl ketone of formula (16).

[0127] In Step G5, the desired compound of formula (I-le) is prepared by deprotecting (desilylating) the compound of formula (17) prepared in Step G4 above with a desilylating agent [e.g. tetrabutylammonium fluoride (TBAF)].

[0128] Each of the reactions in Steps G2 to G5 is well known in the field of synthetic organic chemistry and can be carried out using well known techniques. For example, the reactions in Steps G2 to G5 can be carried out according to the method described in detail in Brian L. Bray et al., J. Org. Chem., 55, 6317-6318 (1990), the contents of which are incorporated herein by reference thereto.

Step G6

[0129] In this Step, an unsaturated heterocycl compound of formula (18) is prepared by subjecting the hydroxy-heterocycl compound of formula (17) prepared in Step G4 above to a dehydration reaction. Dehydration reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques. For example, the dehydration reaction can typically be carried out in the presence of an acid catalyst such as sulfuric acid, a solid catalyst such as alumina or a halogenating agent such as thionyl chloride [e.g. in accordance with the methods described in detail in G. H. Coleman & H. F. Johnstone, Org. Synth., 1, 183 (1941), R. L. Sawyer & D. W. Andrus, Org. Synth., III, 276 (1955) and J. S. Lomas et al., Tetrahedron Lett., 599 (1971), the contents of which are incorporated herein by reference thereto]. Alternatively, the dehydration reaction of this step can be accomplished by a reaction using a trialkylsilane, such as triethylsilane, tripropylsilane or tributylsilane, and trifluoroacetic acid, for example in accordance with the method described in Francis A. Carey & Henry S. Tremper, J. Am. Chem. Soc., 91, 2967-2972 (1969), the contents of which are incorporated herein by reference thereto.

Step G7

[0130] In this Step, the desired compound of formula (I-1f) of the present invention is prepared by removing the protecting group (the silyl group) from the unsaturated heterocycl compound of formula (18) prepared in Step G6 above in a similar manner to that performed in Step G5 above.

Step G8

[0131] In this Step, the desired compound of formula (I-1g) of the present invention is prepared by reducing the double bond in the cyclic group B of the compound of formula (I-1f) prepared in Step G7 above. This is carried out in a similar manner to the procedure described in Step F9 above.

Step G9

[0132] In this Step, a hydroxy compound of formula (20) is prepared by reaction between the bromopyrrole compound of formula (15) prepared in Step G3 above and an aldehyde compound of formula (19). This is carried out in a procedure similar to that described in Step G4 above.

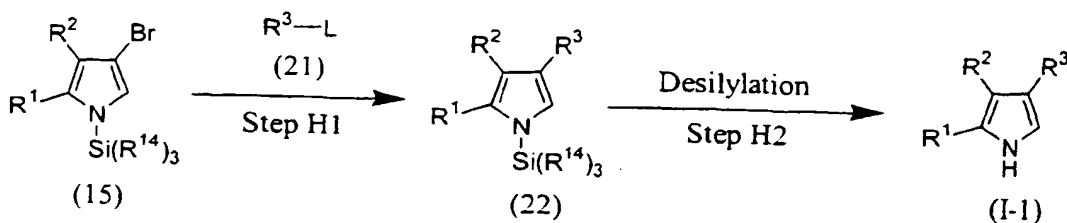
Step G10

[0133] In this Step, the desired compound of formula (I-lh) of the present invention is prepared by removing the protecting group (silyl group) from the hydroxy compound of formula (20) prepared in Step G9 above. This is carried out in a similar manner to that described in Step G5 above.

[0134] Where k in the compound of formula (I-lh) above is an integer of from 1 to 5, the hydroxyalkane moiety can be converted into the corresponding alkenylene moiety by subjecting such a compound to a dehydration reaction in a similar manner to that described in Step G6 above.

Method H

[0135] Generally, the compounds of formula (I) of the present invention can be prepared by introducing the R³ group into a pyrrole compound already substituted on the pyrrole ring with the R¹ group and R² group. Compounds of formula (I-1) can be prepared, for example, according to the Method H, as shown in Reaction Scheme H below.

Reaction Scheme H

[0136] In the above formulae, R¹, R², R³ and R¹⁴ are as defined above, and L represents a leaving group.

[0137] The leaving group L is a group which is capable of leaving as a nucleophilic residue. Examples include halogen atoms such as fluorine, chlorine, bromine and iodine, trihalogenomethoxy groups such as trichloromethoxy, lower alkanesulfonyloxy groups such as methanesulfonyloxy and ethanesulfonyloxy groups, lower halogeno alkane sulfonyloxy groups such as trifluoromethanesulfonyloxy and pentafluoroethanesulfonyloxy groups, and arylsulfonyloxy groups such as benzenesulfonyloxy, p-toluenesulfonyloxy and p-nitrobenzenesulfonyloxy groups. Of these, halogen atoms are preferred, and bromine atoms are particularly preferred.

Step H1

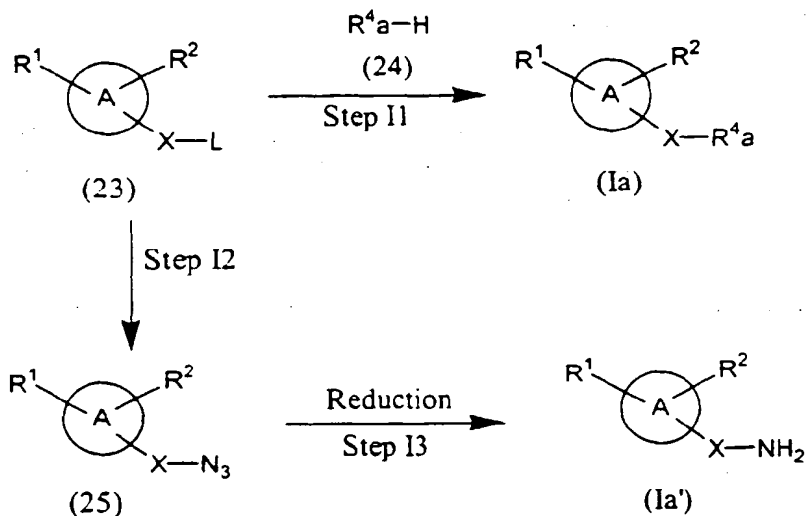
[0138] In this Step, a compound of formula (22) is prepared by lithiating the bromopyrrole compound of formula (15) prepared in Step G3 above using a procedure similar to that described in Step G4 above and then reacting the lithiated intermediate with a compound of formula (21). Substitution reactions of a lithiated pyrrole intermediate of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the methods described in detail in WO 99/01449, the contents of which are incorporated herein by reference thereto.

Step H2

[0139] In this Step, the desired compound of formula (I-1) is prepared by removing the protecting group (silyl group) from the compound of formula (22) prepared in Step H1 above according to a procedure similar to that used in Step G5 above.

Method I

[0140] In this method, compounds of formula (I) of the present invention wherein substituent R^3 is a group of formula $-X-R^4_a$, wherein R^4_a is a heterocyclyl group which may be optionally substituted with at least one group selected from Substituent group α and Substituent group δ and which is substituted with at least one group selected from Substituent group β and Substituent group γ , as defined above, a heterocyclyl group having at least one nitrogen atom, which may be optionally substituted with at least one group selected from Substituent group α and Substituent group δ , as defined above, or a group of formula $-NR^aR^b$, wherein R^a and R^b are as defined above, in which said group is bonded to the group X through the nitrogen atom in said group R^4_a , can be prepared as shown in Reaction Scheme I below.

Reaction Scheme I

[0141] In the above formulae, A, R^1 , R^2 , R^4_a , L and X are as defined above.

Step I1

[0142] In this Step, the desired compound of formula (Ia) of the present invention is prepared by reacting a compound of formula (23) with an amine compound of formula (24) so as to replace the group L with the group R^4_a . Nucleophilic substitution reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques. This reaction is usually carried out in a solvent in the presence or absence of a base.

[0143] There is no particular restriction on the solvent used, provided that it has no adverse effect on the reaction or the reagents involved and that it can dissolve the reagents, at least to some extent. Examples of suitable solvents which can be used include: alcohols such as methanol, ethanol, propanol and isopropanol; ethers such as diethyl ether, diisopropyl ether, tetrahydrofuran and dioxane; aprotic polar solvents such as dimethylformamide, dimethylacetamide and dimethyl sulfoxide; nitriles such as acetonitrile; esters such as methyl acetate and ethyl acetate; aromatic hydrocarbons such as benzene, toluene and xylene; and aliphatic hydrocarbons such as pentane, hexane and heptane.

[0144] Examples of the base which can be used include: alkali metal alkoxides such as sodium methoxide, sodium ethoxide and potassium t-butoxide; alkali metal hydrides such as sodium hydride and lithium hydride; alkali metal hydroxides such as sodium hydroxide and potassium hydroxide; alkali metal carbonates such as sodium carbonate and potassium carbonate; and amines such as triethylamine, tributylamine, pyridine, picoline and 1,8-diazabicyclo [5.4.0]-7-undecene.

[0145] A compound of formula (Ia') of the present invention, which is a compound of formula (Ia) wherein R^4_a is an amino group ($-NH_2$), can also be prepared via an azide compound (25) according to Steps I2 and I3 below.

Step I2

[0146] In this Step, an azide compound of formula (25) is prepared by reacting a compound of formula (23) with sodium azide in the presence of a solvent.

[0147] There is no particular restriction on the solvent used, provided that it has no adverse effect on the reaction or the reagents involved and that it can dissolve the reagents, at least to some extent. Examples of suitable solvents which can be used include: halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride and dichloroethane; ethers such as diethyl ether, diisopropyl ether, tetrahydrofuran and dioxane; aprotic polar solvents such as dimethylformamide, dimethylacetamide and dimethylsulfoxide; nitriles such as acetonitrile; esters such as methyl acetate and ethyl acetate; aromatic hydrocarbons such as benzene, toluene and xylene; aliphatic hydrocarbons such as pentane, hexane and heptane; and mixtures thereof.

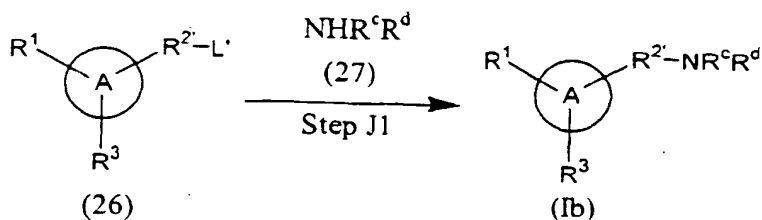
Step I3

[0148] In this Step, the desired compound of formula (Ia') of the present invention is prepared by reducing the azide compound of formula (25) prepared in Step I2 above. Reduction reactions of this type are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the catalytic reduction method described in detail in E. J. Corey et al., Synthesis, 590 (1975), the contents of which are incorporated herein by reference thereto.

Method J

[0149] In this method, compounds of formula (Ib) of the present invention wherein R^2 is a heteroaryl group having at least one nitrogen atom and is substituted with a group of formula $-NR^cR^d$, wherein R^c and R^d are as defined above can be prepared, as shown in Reaction Scheme J below.

Reaction Scheme J



[0150] In the above formulae, the cyclic group A, R^1 , R^3 , R^c and R^d are as defined above, R^2 is a heteroaryl group having at least one nitrogen atom as defined above in the definition of substituent R^2 and L' is a leaving group. Thus, the group $-R^2-L'$ is, for example, 2-methanesulfonylpyrimidin-4-yl, 2-methanesulfonylpyridin-4-yl, etc.

[0151] The leaving group L' is a similar group to the leaving groups defined and exemplified above in the definition of L or it is a lower alkylsulfonyl group as defined and exemplified above, such as a methanesulfonyl, ethanesulfonyl, propanesulfonyl or butanesulfonyl group. The group L' is preferably a lower alkylsulfonyl group, and more preferably a methanesulfonyl group.

Step J1

[0152] In this Step, the desired compound of formula (Ib) of the present invention is prepared by reacting a compound of formula (26) with an amine compound of formula (27) to replace the leaving group with a group of formula $-NR^cR^d$. This step is carried out using a procedure similar to that used in Step I1 above.

Method K

[0153] In this method, compounds of formula (I) of the present invention wherein R^3 is a group containing a pyridine ring can be prepared by subjecting a compound of formula (I) in which R^3 is a tetrahydropyridine group to a dehydrogenation reaction to convert the tetrahydropyridine group to a pyridine group, this dehydrogenation being as shown in Reaction Scheme K below.

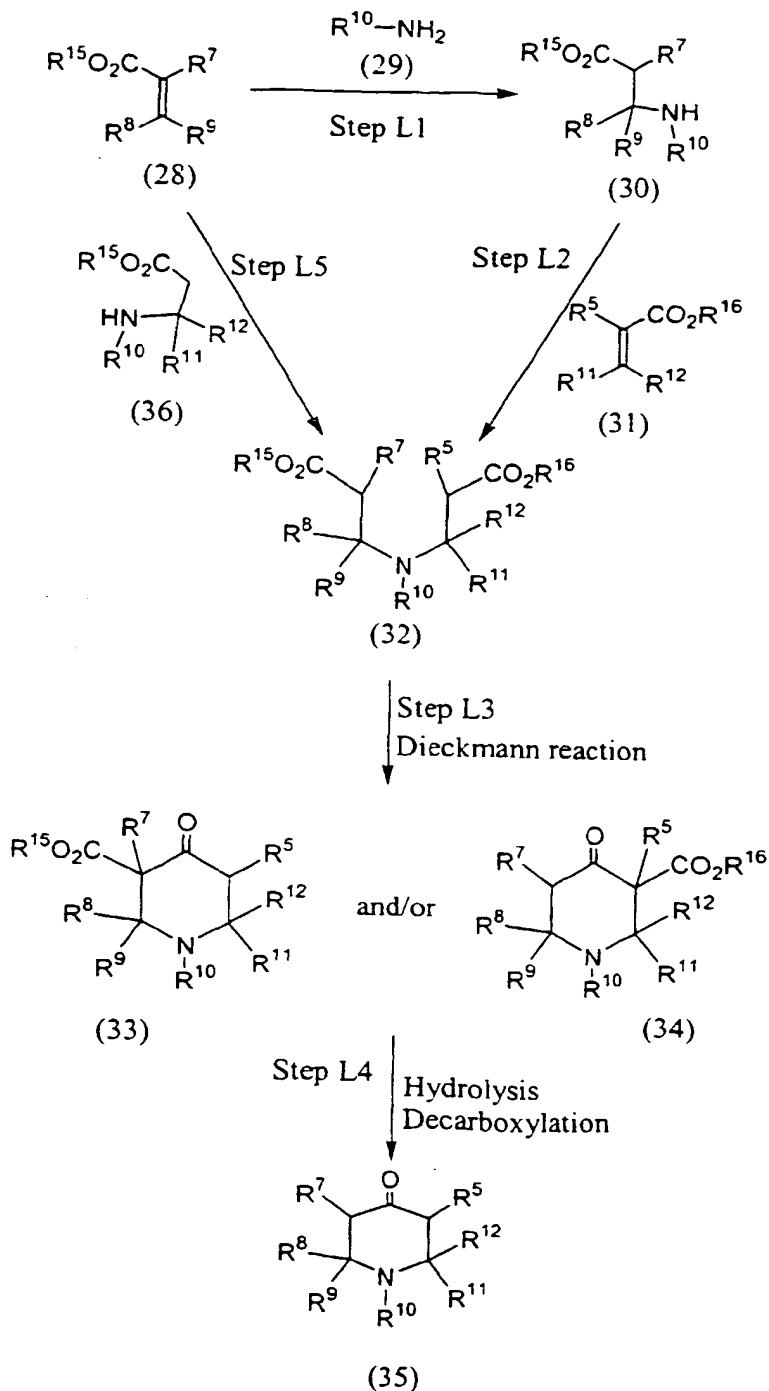
Reaction Scheme K

[0154] Dehydrogenation reactions of the type depicted in Step K1. above are well known in the field of synthetic organic chemistry and can be carried out using well known techniques, for example according to the method described in detail in G. Wieslaw et al., J. Med. Chem., 28, 311-317 (1985), the contents of which are incorporated herein by reference thereto using a palladium-carbon catalyst.

[0155] The compounds which can be used as starting materials in Method A to Method J above, that is the compounds of formulae (1a), (1b), (1c), (1d), (1e), (2a), (2b), (2c), (2d), (2e), (4), (5), (6a), (16), (19), (21), (24) and (27), are either known compounds themselves or are compounds which are easily obtainable by treating known compounds according to known methods (for example, according to the methods described in WO 97/5677, the contents of which are incorporated herein by reference thereto). The starting materials of formulae (23) and (26) in Methods I and J respectively can be easily synthesized from known compounds by carrying out similar reactions to those described in Method A to Method E above.

Method L

[0156] In this method, a 4-oxopiperidine compound of formula (35), which is a starting material falling within the scope of the compounds of formula (16) above, is prepared as shown in the following Reaction Scheme L.

Reaction Scheme L

[0157] In the above formulae, R^5 , R^7 , R^8 , R^9 , R^{10} , R^{11} and R^{12} are as defined above, and R^{15} and R^{16} are the same or different and each is a lower alkyl group as defined above or an aralkyl group as defined above.

Steps L1 to L5

[0158] In step L1, an amino ester compound of formula (30) is prepared by an addition reaction between an unsatu-

rated ester compound of formula (23) and an amine compound of formula (29).

[0159] In Step L2, a diester compound of formula (32) is obtained by a further addition reaction between an unsaturated ester compound of formula (31) and the amino ester compound of formula (30) prepared in Step L1 above.

[0160] In Step L3, the compound of formula (32) prepared in Step L2 above is subjected to Dieckmann reaction to give a keto ester compound of formula (33) and/or formula (34).

[0161] In Step L4, the compound of formula (33) and/or the compound of formula (34) prepared in Step L3 above is subjected to consecutive hydrolysis and decarboxylation reactions to prepare the desired 4-oxopiperidine compound of formula (35).

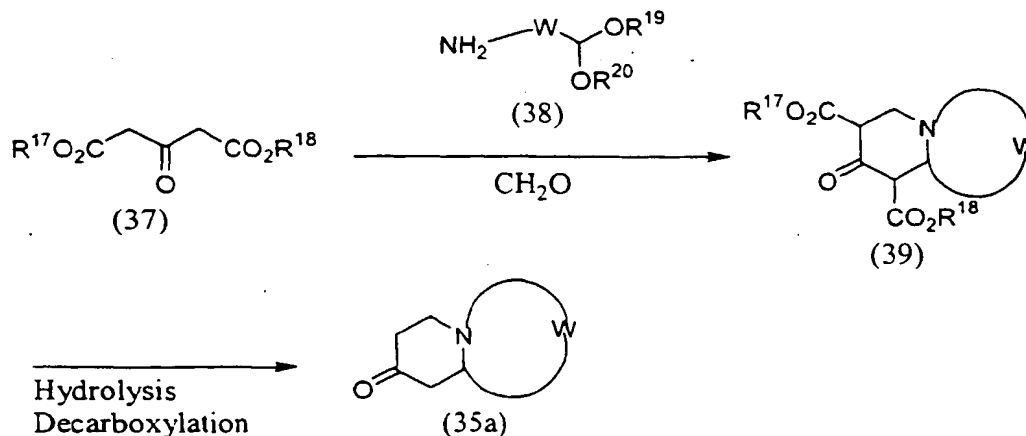
[0162] Step L5 is an alternative process for the preparation of the diester compound of formula (32) and it involves reaction of the unsaturated ester compound of formula (28) with an amino ester compound of formula (36). This is carried out in a similar manner to that described in Step L2 above. Step L5 is preferably employed where the compound of formula (32) to be prepared is one wherein R^{10} and R^{11} together form an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above.

[0163] Each of the reactions in Steps L1 to L5 is well known in the field of synthetic organic chemistry and can be carried out using well known techniques. For example, the reactions in Steps L1 to L5 can be carried out according to the method described in detail in U. M. Teotino, J. Org. Chem., 27, 1406 (1962), the contents of which are incorporated herein by reference thereto.

Method M

[0164] In this method, a compound of formula (35a), which is a 4-oxopiperidine compound of formula (35) above wherein R^{10} and R^{11} together form an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above and each of R^5 , R^7 to R^9 and R^{12} is a hydrogen atom can be prepared as shown in Reaction Scheme M below.

Reaction Scheme M



[0165] In the above formulae, R^{17} and R^{18} may be the same or different and each represents a lower alkyl group as defined and exemplified above or an aralkyl group as defined and exemplified above, R^{19} and R^{20} may be the same or different and each represents a lower alkyl group as defined and exemplified above or R^{19} and R^{20} together represent a lower alkylene group as defined and exemplified above, and W represents an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α defined above, Substituent group β defined above, Substituent group γ defined above and Substituent group δ defined above.

[0166] The two steps of Method M can be carried out according to the methods described in detail in O. Pollet et al., Heterocycles, 43, 1391 (1996) and Anet et al., Austral. J. Scient. Res., <A3>, 635-640 (1950), the contents of which are incorporated herein by reference thereto.

[0167] After completion of each of the reactions described in the steps of Method A to Method M above, the desired compound may be isolated from the reaction mixture in a conventional manner. For example, it can be obtained by neutralizing the reaction mixture as needed, removing insoluble matters by filtration, if any are present, adding organic solvents which are not miscible with each other, such as water and ethyl acetate, washing with water or the like, separating the organic layer containing the desired compound, drying it over anhydrous magnesium sulfate or the like and then distilling off the solvent.

[0168] If necessary, the desired compound thus obtained can be isolated and purified by using a conventional method such as recrystallization or reprecipitation or by a chromatographic method. Examples of chromatography include adsorption column chromatography using a carrier such as silica gel, alumina or magnesium-silica gel type Florisil, chromatography using a synthetic adsorbent, for example, partition column chromatography using a carrier such as Sephadex LH-20 (product of Pharmacia), Amberlite XAD-11 (product of Rohm & Haas) or Diaion HP-20 (product of Mitsubishi Chemical), ion exchange chromatography and normal-phase-reverse-phase column chromatography (high-performance liquid chromatography) using a silica gel or alkylated silica gel. If necessary, two or more of these techniques can be used in combination to isolate and purify the desired compound.

[0169] The heteroaryl-substituted pyrrole derivatives of the present invention exhibit excellent inhibitory activity against the production of inflammatory cytokines. Consequently, they are effective as a medicament, particularly as an agent for the prophylaxis or treatment of diseases mediated by inflammatory cytokines. Examples of such a medicament include analgesics, anti-inflammatory drugs and virucides, and agents for the prophylaxis and treatment of chronic rheumatoid arthritis, osteoarthritis, allergic diseases, asthma, septicemia, psoriasis, osteoporosis, autoimmune diseases (e.g. systemic lupus erythematosus, ulcerative colitis and Crohn's disease), diabetes, nephritis, hepatitis, arteriosclerosis, cancer, ischemic heart disease and Alzheimer's disease. Of these, the compounds of the present invention are particularly useful as analgesics and anti-inflammatory drugs and as agents for the prophylaxis and treatment of chronic rheumatism, osteoarthritis, allergosis, sepsis, psoriasis, osteoporosis, ulcerative colitis, diabetes, hepatitis and arteriosclerosis.

[0170] The compounds of formula (I) and pharmacologically acceptable salts, esters and other derivatives thereof according to the present invention can be administered by a number of different routes. Examples of these administration routes include oral administration in the form of tablets, capsules, granules, powders or syrups and parenteral administration in the form of injections or suppositories. Such formulations can be prepared in a known manner by using additives such as an excipients, lubricants, binders, disintegrators, stabilizers, corrigents and diluents.

[0171] Examples of suitable excipients include: organic excipients, examples of which include sugar derivatives such as lactose, sucrose, dextrose, mannitol and sorbitol, starch derivatives such as corn starch, potato starch, α -starch, dextrin and carboxymethyl starch, cellulose derivatives such as crystalline cellulose, low-substituted hydroxypropylcellulose, hydroxypropylmethylcellulose, carboxymethylcellulose, calcium carboxymethylcellulose and sodium internally-crosslinked carboxymethylcellulose, gum arabic, dextran and pullulan; and inorganic excipients, examples of which include silicate derivatives such as soft silicic acid anhydride, synthetic aluminum silicate and magnesium aluminometasilicate, phosphates such as calcium phosphate, carbonates such as calcium carbonate, and sulfates such as calcium sulfate.

[0172] Examples of suitable lubricants include: stearic acid; metal salts of stearic acid such as calcium stearate and magnesium stearate; talc; colloidal silica; waxes such as bee gum and spermaceti; boric acid; adipic acid; sulfates such as sodium sulfate; glycol; fumaric acid; sodium benzoate; DL-leucine; sodium salts of an aliphatic acid; lauryl sulfates such as sodium lauryl sulfate and magnesium lauryl sulfate; silicic acid derivatives such as silicic anhydride and silicic acid hydrate; and starch derivatives exemplified above as examples of suitable excipients.

[0173] Examples of suitable binders include polyvinylpyrrolidone, Macrogol and compounds similar to those exemplified above as suitable excipients.

[0174] Examples of suitable disintegrators include compounds similar to those exemplified above as suitable excipients and chemically modified starch or cellulose derivatives such as sodium cross carmellose, sodium carboxymethyl starch and crosslinked polyvinylpyrrolidone.

[0175] Examples of suitable stabilizers include: paraoxybenzoate esters such as methylparaben and propylparaben; alcohols such as chlorobutanol, benzyl alcohol and phenylethyl alcohol; benzalkonium chloride; phenol derivatives such as phenol and cresol; thimerosal; dehydroacetic acid; and sorbic acid. Examples of suitable corrigents include sweeteners, acidifiers and flavors commonly employed for this purpose.

[0176] The dose of the compound of formula (I) or a pharmacologically acceptable salt, ester or other derivative thereof according to the present invention will vary depending on a variety of factors including the condition to be treated, the age of the patient and the administration route. When administered orally, it is administered to an adult in an amount of 0.1 mg (preferably 0.5 mg) a day as a lower limit and 2000 mg (preferably 500 mg) a day as an upper limit. It can be administered in from one to several portions depending on the condition of the patient. When administered intravenously, it is administered to an adult in an amount of 0.01 mg (preferably 0.05 mg) a day as a lower limit and 200 mg (preferably 50 mg) a day as an upper limit. It can be administered in from one to several portions depending

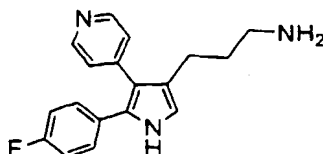
on the condition of the patient.

[0177] The following examples, preparative examples, formulation examples and test examples are intended to further illustrate the present invention and are not intended to limit the scope of this invention in any way.

Example 1

4-(3-Aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-97)

[0178]



1(i) 4-Ethoxycarbonyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0179] 36 ml (54.7 mmol) of a 1.53 N solution of butyllithium in hexane were added to 240 ml of tetrahydrofuran. A solution of 15.90 g (54.7 mmol) of α -(p-toluene-sulfonyl)-4-fluorobenzylisocyanide in 120 ml of tetrahydrofuran was then added to the resulting solution at -45°C , followed by stirring of the resulting mixture for 10 minutes at the same temperature. At the end of this time, 25.00 g (273 mmol) of 95% lithium bromide were added, the resulting mixture was stirred for 30 minutes and then a solution of 8.73 g (49.2 mmol) of ethyl 3-(4-pyridyl)acrylate in 120 ml of tetrahydrofuran was added. The resulting mixture was stirred at the same temperature for 1 hour and then the cooling bath was removed and the mixture was stirred at room temperature for a further 1 hour. At the end of this time, 500 ml of water were added and the resulting mixture was extracted with ethyl acetate. The organic extract was washed with water and then dried over anhydrous sodium sulfate, after which it was concentrated by evaporation under reduced pressure to afford a solid. The solid was washed with diethyl ether to give 13.61 g (yield 89%) of the title compound as a pale yellow powder.

[0180] ^1H -Nuclear magnetic resonance spectrum (500 MHz, CDCl_3) δ ppm:

8.34 (1H, broad singlet);
8.51 (2H, doublet, $J=7$ Hz);
7.58 (1H, doublet, $J=3$ Hz);
7.21 (2H, doublet, $J=6$ Hz);
7.11 (2H, doublet of doublets, $J=9$ Hz, 5 Hz);
6.97 (2H, triplet, $J=9$ Hz);
4.18 (2H, quartet, $J=7$ Hz);
1.20 (3H, triplet, $J=7$ Hz).

1(ii) 2-(4-Fluorophenyl)-4-hydroxymethyl-3-(pyridin-4-yl)-1H-pyrrole

[0181] 121.4 ml (121.4 mmol) of a 1M solution of diisobutylaluminum hydride in toluene were slowly added dropwise to a solution of 12.56 g (40.47 mmol) of 4-ethoxycarbonyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [which was obtained in step (i) above] in 630 ml of tetrahydrofuran with ice-cooling. After completing the addition, the cooling bath was removed and the mixture was stirred at room temperature for 5 hours. At the end of this time, a saturated aqueous solution of ammonium chloride was added to the reaction mixture and the resulting mixture was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure to give 10.30 g (yield 99%) of the title compound as a pale orange powder.

[0182] ^1H -Nuclear magnetic resonance spectrum (500 MHz, CDCl_3) δ ppm:

10.52 (1H, broad singlet);
8.46 (2H, doublet, $J=6$ Hz);
7.31 (2H, doublet, $J=6$ Hz);
7.26 (2H, doublet of doublets, $J=9$ Hz, 5 Hz);
6.98 (2H, triplet, $J=9$ Hz);
6.95 (1H, doublet, $J=3$ Hz);

4.50 (2H, singlet);
3.49 (1H, broad singlet).

1(iii) 2-(4-Fluorophenyl)-4-formyl-3-(pyridin-4-yl)-1-H-pyrrole

[0183] 118.00 g (440 mmol) of manganese dioxide were added to a solution of 10.73 g (40 mmol) of 2-(4-fluorophenyl)-4-hydroxymethyl-3-(pyridin-4-yl)-1-H-pyrrole [prepared as described in step (ii) above] in 120 ml of dimethyl sulfoxide, after which the resulting mixture was stirred at 50°C overnight. At the end of this time, the reaction mixture was filtered, water was added to the filtrate, and this was then extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure to afford a solid which was washed with diethyl ether to give 7.35 g (yield 69%) of the title compound as a brown powder.

[0184] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃-CD₃OD) δ ppm:

9.76 (1H, singlet);
8.48 (2H, doublet, J=6 Hz);
7.59 (1H, singlet);
7.26 (2H, doublet, J=6 Hz);
7.19 (2H, doublet of doublets, J=9 Hz, 6 Hz);
7.01 (2H, triplet, J=9 Hz).

1(iv) 3-[2-(4-Fluorophenyl)-3-(pyridin-4-yl)-1-H-pyrrol-4-yl]acrylonitrile

[0185] 2.92 ml (18.02 mmol) of diethylphosphonoacetonitrile were added to 16 ml of tetrahydrofuran and the resulting solution was added to a suspension of 786 mg (18.02 mmol) of 55% sodium hydride in 60 ml of tetrahydrofuran with ice-cooling, and the resulting mixture was then stirred at room temperature for 1.5 hours. At the end of this time, a suspension of 4.00 g (15.02 mmol) of 2-(4-fluorophenyl)-4-formyl-3-(pyridin-4-yl)-1-H-pyrrole [prepared as described in step (iii) above] in 30 ml of tetrahydrofuran was added to this mixture and the resulting mixture was stirred at room temperature for 1 hour. At the end of this time, water was added to the reaction mixture and the resulting mixture was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure to afford a solid which was washed with diethyl ether to give 2.94 g (yield 68%) of the title compound as a pale brown powder.

[0186] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃-DMSO-d₆) δ ppm:

11.30 (1H, broad singlet);
8.57 (2H, doublet, J=6 Hz);
7.24 (1H, doublet, J=4 Hz);
7.21 (1H, doublet, J=17 Hz);
7.19 (2H, doublet of doublets, J=9 Hz, 5 Hz);
7.13 (2H, doublet, J=6 Hz);
6.97 (2H, doublet, J=9 Hz);
5.39 (1H, doublet, J=17 Hz).

1(v) 4-(2-Cyanoethyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-H-pyrrole

[0187] 2.94 g (10.16 mmol) of 3-[2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-H-pyrrol-4-yl]acrylonitrile [prepared as described in step (iv) above] were dissolved in a mixture of 20 ml of tetrahydrofuran and 20 ml of methanol, after which 2.94 g of 10% palladium on carbon were added to the solution. The mixture was stirred under a hydrogen atmosphere at room temperature for 8 hours. At the end of this time, the reaction mixture was filtered and the filtrate was concentrated under reduced pressure. The resulting residue was washed with a mixture of 30 ml of diethyl ether and 30 ml of ethanol to give 1.80 g (yield 61%) of the title compound as a pale orange powder.

[0188] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.55 (2H, doublet, J=6 Hz);
8.27 (1H, broad singlet);
7.13 (2H, doublet of doublets, J=9 Hz, 5 Hz);
7.12 (2H, doublet, J=6 Hz);
6.98 (2H, triplet, J=9 Hz);
6.39 (1H, doublet, J=3 Hz);
2.38 (2H, triplet, J=7 Hz);
2.43 (2H, triplet, J=7 Hz).

1(vi) 4-(3-Aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0189] 1.80 g (6.18 mmol) of 4-(2-cyanoethyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step (v) above] were added to 100 ml of tetrahydrofuran and the resulting solution was slowly added dropwise with stirring at room temperature to a suspension of 469 mg (12.36 mmol) of lithium aluminum hydride in 150 ml of tetrahydrofuran. This mixture was stirred at 60°C for 30 minutes. At the end of this time, the reaction mixture was cooled to room temperature, 25 ml of water and 0.5 ml of a 15% aqueous solution of sodium hydroxide were slowly added to the reaction mixture and the resulting mixture was then filtered. The filtrate was concentrated under reduced pressure, after which the residue was dehydrated by azeotropic distillation under reduced pressure using toluene. This azeotropic distillation was carried out three times to give 1.71 g (yield 94%) of the title compound as a white powder.

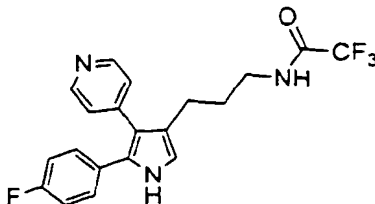
Melting point: 198 - 199°C

[0190] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.16 (1H, broad singlet);
 8.46 (2H, doublet, J=6 Hz);
 7.18 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 7.13 (2H, doublet, J=6 Hz);
 7.11 (2H, triplet, J=9 Hz);
 6.74 (1H, doublet, J=3 Hz);
 2.49 (2H, triplet, J=7 Hz);
 2.41 (2H, triplet, J=6 Hz);
 1.48 (2H, quintet, J=8 Hz).

Example 22-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(3-trifluoroacetylaminopropyl)-1H-pyrrole (Compound No. 1-2533)

[0191]



[0192] 3.3 ml (23.2 mmol) of trifluoroacetic anhydride were added to a solution of 685 mg (2.32 mmol) of 4-(3-aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (which was prepared as described in Example 1 above) in 70 ml of tetrahydrofuran. The resulting mixture was stirred at room temperature for 30 minutes. At the end of this time, 200 ml of a saturated aqueous solution of sodium hydrogencarbonate were added to the reaction mixture, and then this was extracted with ethyl acetate. The organic extract was washed with water and concentrated by evaporation under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using a 1:1 by volume mixture of hexane and ethyl acetate as the eluant to give 280 mg (yield 31%) of the title compound as a white powder.

[0193] Melting point: 229 - 230°C

[0194] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

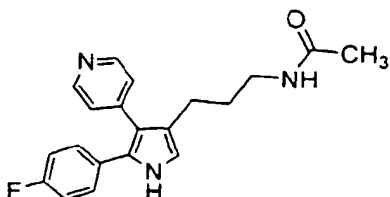
10.09 (1H, broad singlet);
 8.49 (2H, doublet, J=6 Hz);
 7.93 (1H, broad singlet);
 7.18 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.13 (2H, doublet, J=6 Hz);
 6.94 (2H, triplet, J=9 Hz);
 6.74 (1H, doublet, J=3 Hz);
 3.29 (2H, doublet of doublets, J=13 Hz, 7 Hz);

2.56 (2H, triplet, $J=8$ Hz);
1.74 (2H, quintet, $J=8$ Hz).

Example 3

4-(3-Acetylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2526)

[0195]



[0196] In a similar manner to that described in Example 2 above, a reaction was carried out using acetic anhydride instead of trifluoroacetic anhydride to give the title compound (yield 99%) as a pale yellow powder.

Melting point: 218 -220°C

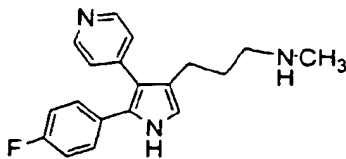
[0197] ^1H -Nuclear magnetic resonance spectrum (400 MHz, DMSO-d_6) δ ppm:

11.19 (1H, broad singlet);
8.46 (2H, doublet, $J=6$ Hz);
7.77 (1H, broad triplet, $J=5$ Hz);
7.18 (2H, doublet of doublets, $J=9$ Hz, 5 Hz);
7.15-7.08 (4H, multiplet);
6.77 (1H, d, 2 Hz);
2.99 (2H, doublet of doublets, $J=13$ Hz, 7 Hz);
2.39 (2H, triplet, $J=8$ Hz);
1.75 (3H, singlet);
1.53 (2H, quintet, $J=7$ Hz).

Example 4

2-(4-Fluorophenyl)-4-(3-methylaminopropyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-101)

[0198]



4(i) 4-[3-(t-Butoxycarbonylamino)propyl]-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0199] In a similar manner to that described in Example 2 above, a reaction was carried out using di-*t*-butyl dicarbonate instead of trifluoroacetic anhydride and using methanol instead of tetrahydrofuran to give the title compound (yield 68%) as a white powder.

[0200] ^1H -Nuclear magnetic resonance spectrum (400 MHz, CDCl_3) δ ppm:

8.51 (2H, doublet, $J=6$ Hz);
8.26 (1H, broad singlet);

7.15-7.12 (4H, multiplet);
 6.96 (2H, triplet, J=9 Hz);
 6.75 (1H, doublet, J=3 Hz);
 4.45 (1H, broad singlet);
 3.12-3.09 (2H, multiplet);
 2.54 (2H, triplet, J=8 Hz);
 1.67-1.63 (2H, multiplet);
 1.43 (9H, singlet).

4(ii) 2-(4-Fluorophenyl)-4-(3-methylaminopropyl)-3-(pyridin-4-yl)-1H-pyrrole

[0201] 400 mg (1.01 mmol) of 4-[3-(t-butoxycarbonylamino)propyl]-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 4(i) above] were added to a suspension of 154 mg of lithium aluminum hydride in 8 ml of tetrahydrofuran. The resulting mixture was heated under reflux for 2 hours. At the end of this time, the reaction mixture was worked up and then purified in the same manner as that described in Example 1(vi) above to give 300 mg (yield 95%) of the title compound as a white powder.

[0202] Melting point: 193-198°C

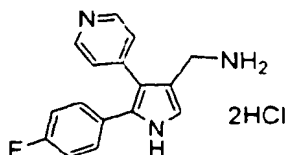
[0203] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.51 (2H, doublet, J=6 Hz);
 8.34 (1H, broad singlet);
 7.17-7.12 (4H, multiplet);
 6.96 (2H, triplet, J=9 Hz);
 6.74 (1H, doublet, J=3 Hz);
 2.57 (2H, triplet, J=7 Hz);
 2.54 (2H, triplet, J=7 Hz);
 2.38 (3H, singlet);
 1.69 (2H, quintet, J=7 Hz).

Example 5

4-Aminomethyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole dihydrochloride (Dihydrochloride of Compound No. 1-95)

[0204]



5(i) 5-(4-Fluorophenyl)-4-(pyridin-4-yl)-1H-pyrrole-3-carboxaldehyde O-methyloxime

[0205] 0.83 g (9.58 mmol) of O-methylhydroxylamine hydrochloride and 1.56 ml (11.2 mmol) of triethylamine were added to a solution of 0.85 mg (3.2 mmol) of 2-(4-fluorophenyl)-4-formyl-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 1(iii) above] in 17 ml of methanol. The resulting mixture was heated under reflux for 30 minutes, after which water was added to the reaction mixture which was then extracted with ethyl acetate. The organic extract was washed with water and concentrated by evaporation under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using ethyl acetate as the eluant to afford 728 mg (yield 77%) of the title compound as a pale yellow powder.

[0206] Melting point: 242 - 245°C

[0207] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.88 (1H, broad singlet);
 8.54 (0.2H, doublet, J=6 Hz);
 8.49 (1.8H, doublet, J=6 Hz);
 7.88 (0.9H, singlet);
 7.76 (0.1 H, singlet);

7.34 (0.9H, singlet);
 7.25-7.12 (6H, multiplet);
 6.98 (0.1H, singlet);
 3.92 (0.4H, singlet);
 3.71 (2.6H, singlet).

5(ii) 4-Aminomethyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole dihydrochloride

[0208] 10 ml of 10% hydrochloric acid and 0.50 g of 10% palladium on carbon were added to a solution of 497 mg (1.68 mmol) of 5-(4-fluorophenyl)-4-(pyridin-4-yl)-1H-pyrrole-3-carboxaldehyde *O*-methyloxime [which was prepared as described in step 5(i) above] in 10 ml of methanol. The mixture was stirred under a hydrogen atmosphere at room temperature for 2 hours. At the end of this time, the reaction mixture was filtered and the filtrate thus obtained was concentrated by evaporation under reduced pressure. The residue was washed with diisopropylether and then with diethylether to give 391 mg (yield 68%) of the title compound as a yellowish orange powder.

[0209] Melting point: 235 -240°C (decomposition)

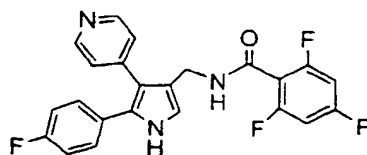
[0210] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

11.85 (1H, broad singlet);
 8.65 (2H, doublet, J=7 Hz);
 7.31 (2H, doublet, J=7 Hz);
 7.32 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.24 (1H, doublet, J=3 Hz);
 7.14 (2H, triplet, J=9 Hz);
 4.22 (2H, singlet).

Example 6

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(2,4,6-trifluorobenzoylaminoethyl)-1H-pyrrole (Compound No. 1-2535)

[0211]



[0212] 265 µl (1.905 mmol) of triethylamine and 96 mg (0.493 mmol) of 2,4,6-trifluorobenzoyl chloride were added to a suspension of 162 mg (0.476 mmol) of 4-aminomethyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole dihydrochloride [prepared as described in Example 5 above] in 3.3 ml of tetrahydrofuran. The resulting mixture was stirred at room temperature for 4 hours. At the end of this time, water was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was washed with diethylether to give 78 mg (yield 38%) of the title compound as a pale yellow powder.

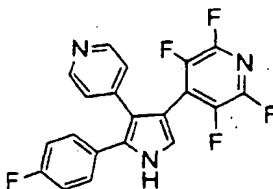
[0213] Melting point: 243 -245°C (decomposition)

[0214] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.41 (1H, broad singlet);
 8.99 (1H, broad triplet, J=5 Hz);
 8.46 (2H, doublet, J=6 Hz);
 7.29-7.19 (4H, multiplet);
 7.19-7.13 (4H, multiplet);
 6.94 (1H, doublet, J=3 Hz);
 4.27 (2H, doublet, J=5 Hz).

Example 72-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(2,3,5,6-tetrafluoropyridin-4-yl)-1H-pyrrole (Compound No. 1-2538)

[0215]

7(i) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0216] 15.00 g (48.3 mmol) of 4-ethoxycarbonyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 1(i) above] were dissolved in a mixture of 90 ml of acetic acid, 30 ml of sulfuric acid and 60 ml of water and the resulting solution was stirred at 100°C for 16 hours. At the end of this time, the reaction mixture was cooled to room temperature and then made alkaline with a 10% aqueous solution of sodium hydroxide. The resulting mixture was extracted with ethyl acetate and the organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure to give 11.40 g (yield 99%) of the title compound as a pale red powder.

[0217] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

9.78 (1H, broad singlet);
 8.42 (2H, doublet, J=7 Hz);
 7.37 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.22 (2H, doublet, J=6 Hz);
 7.06 (2H, triplet, J=9 Hz);
 6.90 (1H, triplet, J=3 Hz);
 6.47 (1H, triplet, J=3 Hz).

7(ii) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0218] 31 ml (47.4 mmol) of a 1.57 N solution of butyllithium in hexane were added to a solution of 11.30 g (47.4 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 7(i) above] in 300 ml of tetrahydrofuran at -78°C and the mixture was stirred for 10 minutes. At the end of this time, 13.4 ml (49.8 mmol) of triisopropylsilyl triflate were added to the reaction mixture at the same temperature. After removal of the cooling bath, the mixture was stirred at room temperature for 30 minutes. 200 ml of water and 300 ml of a saturated aqueous solution of sodium hydrogencarbonate were then added to the reaction mixture before extracting with ethyl acetate. The organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure to give 18.70 g (quantitative yield) of the title compound as a reddish purple oil.

[0219] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

8.25 (2H, doublet, J=6 Hz);
 7.39 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 7.28 (2H, triplet, J=9 Hz);
 7.00 (1H, doublet, J=3 Hz);
 6.91 (2H, doublet, J=7 Hz);
 6.71 (1H, doublet, J=3 Hz);
 1.15-1.05 (3H, multiplet);
 0.98 (18H, doublet, J=6 Hz).

7(iii) 4-Bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0220] A suspension of 8.61 g (47.4 mmol) of N-bromosuccinimide in 100 ml of tetrahydrofuran was added to a solution of 18.70 g (47.4 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as de-

scribed in step 7(ii) above] in 300 ml of tetrahydrofuran at -78°C, after which the resulting mixture was stirred at the same temperature for 6 hours. After removal of the cooling bath, the mixture was then stirred at room temperature for a further 1 hour. At the end of this time, 400 ml of hexane was added to the reaction mixture and the insoluble material was removed by filtration. The resulting filtrate was concentrated by evaporation under reduced pressure and the residue thus obtained was purified by chromatography on a silica gel column using a 2:1 by volume mixture of hexane and ethyl acetate as the eluant to give 9.57 g (yield 43%) of the title compound as pale yellow prismatic crystals.

[0221] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

8.36 (2H, doublet, J=6 Hz);
7.34 (2H, doublet of doublets, J=9 Hz, 6 Hz);
7.18 (2H, triplet, J=9 Hz);
7.12 (1H, singlet);
7.04 (2H, doublet, J=6 Hz);
1.16-1.08 (3H, multiplet);
0.99 (16H, doublet, J=6 Hz).

7(iv) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(2,3,5,6-tetrafluoropyridin-4-yl)-1H-pyrrole

[0222] 1.8 ml (2.75 mmol) of 1.54 N solution of t-butyllithium in pentane were added to a solution of 650 mg (1.37 mmol) of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 7(iii) above] in 15 ml of tetrahydrofuran at -78°C and the mixture was stirred for 10 minutes. At the end of this time, 165 µl (1.51 mmol) of pentafluoropyridine were added to the reaction mixture. The cooling bath was then removed and the mixture was stirred for 30 minutes. At the end of this time, 70 ml of water and 50 ml of a saturated aqueous solution of sodium hydrogencarbonate were added to the reaction mixture, and then this was extracted with ethyl acetate. The organic extract was washed with water and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 4:1 by volume mixture of hexane and ethyl acetate as the eluant to give 346 mg (yield 65%) of the title compound as a pale yellow powder.

[0223] Melting point: >300°

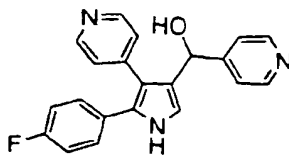
[0224] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.34 (2H, doublet, J=8 Hz);
7.32 (1H, singlet);
7.30 (2H, doublet of doublets, J=11 Hz, 7 Hz);
7.09 (2H, triplet, J=11 Hz);
7.06 (2H, doublet, J=8 Hz).

Example 8

(±)-[5-(4-Fluorophenyl)-4-(pyridin-4-yl)-1H-pyrrol-3-yl]-(pyridin-4-yl)methanol (Compound No. 1-2530)

[0225]



8(i) (±)-[5-(4-Fluorophenyl)-4-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrol-3-yl]-(pyridin-4-yl)methanol

[0226] In a similar manner to that described in Example 7(iv) above, a reaction was carried out using 4-formylpyridine instead of pentafluoropyridine to give the title compound (yield 56%) as a pale yellow oil.

[0227] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.39 (2H, doublet, J=7 Hz);
8.21 (2H, doublet, J=6 Hz);
7.38 (2H, doublet, J=7 Hz);
7.29 (2H, doublet of doublets, J=9 Hz, 6 Hz);
7.17 (2H, doublet, J=6 Hz);

7.03 (2H, triplet, J=9 Hz);
 6.67 (1H, singlet);
 5.32 (1H, singlet);
 1.14-1.05 (3H, multiplet);
 1.01 (9H, doublet, J=7 Hz);
 1.00 (9H, doublet, J=7 Hz).

8(ii) (±)-[5-(4-Fluorophenyl)-4-(pyridin-4-yl)-1H-pyrrol-3-yl]-(pyridin-4-yl)methanol

[0228] 0.59 ml (0.59 mmol) of a 1 M solution of tetrabutylammonium fluoride in tetrahydrofuran were added to a solution of 145 mg (0.295 mmol) of [5-(4-fluorophenyl)-4-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrol-3-yl]-(pyridin-4-yl)methanol [prepared as described in step 8(i) above] in 3 ml of tetrahydrofuran. The resulting mixture was stirred at room temperature for 10 minutes, after which water was added to the reaction mixture and the resulting mixture was acidified with 1 N hydrochloric acid and then extracted with ethyl acetate. The water layer was made alkaline with potassium carbonate and then extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure. The residual oil thus obtained was solidified by the addition of isopropanol to give 83 mg (yield 82%) of the title compound as a white powder.

[0229] Melting point 202 - 205°C

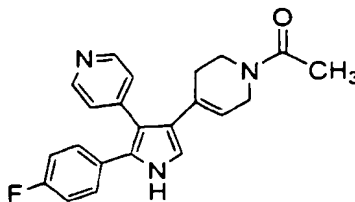
[0230] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.39 (1H, broad singlet);
 8.43 (2H, doublet, J=6 Hz);
 8.41 (2H, doublet, J=6 Hz);
 7.26-7.21 (4H, multiplet);
 7.18 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 7.13 (2H, triplet, J=9 Hz);
 6.55 (1H, doublet, J=3 Hz);
 5.77 (1H, doublet, J=5 Hz);
 5.54 (1H, doublet, J=4 Hz).

Example 9

4-(1-Acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2525)

[0231]



9(i) 4-(1-Benzyl-4-hydroxypiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0232] In a similar manner to that described in Example 7(iv) above, a reaction was carried out using 1-benzylpiperidine-4-one instead of pentafluoropyridine to afford the title compound (yield 61%) as white needle-like crystals.

[0233] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.21 (2H, doublet, J=6 Hz);
 7.35-7.15 (9H, multiplet);
 6.93 (2H, triplet, J=9 Hz);
 6.90 (1H, singlet);
 3.49 (2H, singlet);
 2.58-2.41 (4H, broad multiplet);
 2.00-1.91 (2H, broad multiplet);
 1.83-1.74 (2H, broad multiplet);
 1.20-1.10 (3H, multiplet);

1.06 (1H, doublet, J=7 Hz).

9(ii) 2-(4-Fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

5 [0234] 0.58 g of 10% palladium on carbon were added to a solution of 0.58 g (1.00 mmol) of 4-(1-benzyl-4-hydroxypiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 9(i) above] in 15 ml of methanol. The resulting mixture was stirred under a hydrogen atmosphere at room temperature for 5 hours. At the end of this time, the reaction mixture was filtered and the filtrate was concentrated by evaporation under reduced pressure to give 483 mg (yield 98%) of the title compound as a white amorphous solid.

10 [0235] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:
 8.22 (2H, doublet, J=7 Hz);
 7.27 (2H, doublet, J=6 Hz);
 7.21 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 6.94 (2H, triplet, J=9 Hz);
 15 6.92 (1H, singlet);
 2.98-2.90 (2H, multiplet);
 2.76-2.68 (2H, multiplet);
 1.93-1.84 (2H, multiplet);
 1.77-1.69 (2H, multiplet);
 20 1.20-1.11 (3H, multiplet);
 1.07 (1H, doublet, J=7 Hz).

9(iii) 4-(1-Acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

25 [0236] 1.4 ml (8.70 mmol) of triethylsilane and 2.25 ml (29.0 mmol) of trifluoroacetic acid were added to a solution of 3.58 g (7.25 mmol) of 2-(4-fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 9(ii) above] in 40 ml of dichloromethane, after which the resulting mixture was stirred at room temperature for 5 hours. At the end of this time, a saturated aqueous solution of sodium hydrogencarbonate was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure to afford 3.90 g of dehydrated product as a yellow amorphous solid.

30 [0237] All of the 3.90 g of dehydrated product thus obtained was dissolved in 80 ml of tetrahydrofuran, and then 0.76 ml (8.0 mmol) of acetic anhydride and 2.25 ml (16.0 mmol) of triethylamine were added to the solution. The resulting mixture was stirred at room temperature for 20 minutes, after which 16 ml (16.0 mmol) of a 1 M solution of tetrabutylammonium fluoride in tetrahydrofuran were added before stirring for a further 10 minutes. 150 ml of a saturated aqueous solution of sodium hydrogencarbonate were added to the reaction mixture and then this was extracted with ethyl acetate. After washing the organic extract with water, 1 N hydrochloric acid was added thereto. The water layer was washed with ethyl acetate, made alkaline with potassium carbonate and finally extracted with ethyl acetate. The organic extract thus obtained was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure to give 1.73 g (yield 68%) of the title compound as a pale brown powder.

40 [0238] Melting point: 262 - 264°C (decomposition)

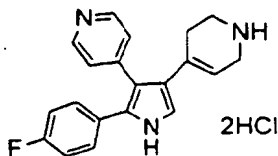
[0239] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.43 (1H, broad singlet);
 8.48-8.44 (2H, multiplet);
 45 7.20-7.09 (6H, multiplet);
 6.96 (0.5H, doublet, J=3 Hz);
 6.94 (0.5H, doublet, J=3 Hz);
 5.32-5.29 (0.5H, broad multiplet);
 5.29-5.25 (0.5H, broad multiplet);
 50 3.92-3.88 (1H, multiplet);
 3.37-3.84 (1H, multiplet);
 3.50 (1H, triplet, J=6 Hz);
 3.46 (1H, triplet, J=6 Hz);
 2.22-2.17 (1H, broad multiplet);
 55 2.16-2.11 (1H, broad multiplet);
 2.00 (1.5H, singlet);
 1.96 (1.5H, singlet).

Example 10

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole dihydrochloride (dihydrochloride of Compound No. 1-143)

[0240]



[0241] 249 mg (0.689 mmol) of 4-(1-acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 9 above] were dissolved in a mixture of 2.5 ml of methanol and 2.5 ml of water, after which 2.43 ml (9.646 mmol) of a 4 N solution of hydrogen chloride in dioxane were added to the solution. The resulting mixture was heated under reflux for 11 hours, after which the reaction mixture was concentrated by evaporation under reduced pressure and the residual solid was washed with hot isopropanol to give 216 mg (yield 80%) of the title compound as a pale yellow powder.

[0242] Melting point: 290 - 295°C (decomposition)

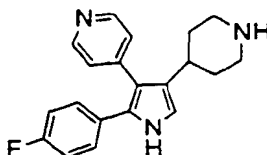
[0243] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

11.63 (1H, broad singlet);
 8.54 (2H, doublet, J=7 Hz);
 7.74 (2H, doublet, J=7 Hz);
 7.29 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.14 (2H, triplet, J=9 Hz);
 7.06 (1H, doublet, J=2 Hz);
 5.51-5.47 (1H, broad multiplet);
 3.72 (2H, doublet, J=3 Hz);
 3.42 (2H, triplet, J=6 Hz);
 2.66-2.55 (2H, broad multiplet).

Example 11

2-(4-Fluorophenyl)-4-(piperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-142)

[0244]



[0245] 120 mg of 10% palladium on carbon were added to a solution of 120 mg (0.306 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole dihydrochloride [prepared as described in Example 10 above] in 10 ml of methanol. The resulting mixture was stirred under a hydrogen atmosphere at room temperature for 1 hour, after which the reaction mixture was filtered. The filtrate thus obtained was concentrated by evaporation under reduced pressure. Water was added to the resulting residue and then this mixture was made alkaline with a 1 N aqueous solution of sodium hydroxide. The resulting precipitate was collected by filtration to give 84 mg (yield 85%) of the title compound as a pale red powder.

[0246] Melting point: 265 - 267°C (decomposition)

[0247] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

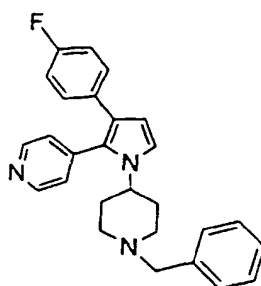
11.17 (1H, broad singlet);

8.48 (2H, doublet, J=6 Hz);
 7.20-7.04 (6H, multiplet);
 6.73 (1H, doublet, J=3 Hz);
 2.38 (2H, doublet, J=12 Hz);
 2.46 (1H, triplet of triplets, J=12 Hz, 4 Hz);
 2.38 (2H, triplet, J=12 Hz);
 1.58 (2H, doublet, J=13 Hz);
 1.35 (2H, double doublet of triplets, J=13 Hz, 12 Hz, 4 Hz).

Example 12

1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-147)

[0248]



12(i) 1-Benzyl-4-(pyridin-4-yl)methyleaminopiperidine

[0249] 8.57 ml (42.01 mmol) of 4-amino-1-benzylpiperidine were added to a solution of 3.95 ml (42.01 mmol) of 4-formylpyridine in 5 ml of ethanol, and the resulting mixture was then heated under reflux for 1 hour. At the end of this time, the solvent was distilled off from the reaction mixture by evaporation under reduced pressure to afford 11.78 g (quantitative yield) of the title compound as a pale brown powder.

[0250] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.68 (2H, doublet, J=6 Hz);
 8.30 (1H, singlet);
 7.60 (2H, doublet, J=6 Hz);
 7.38-7.22 (5H, multiplet);
 3.56 (2H, singlet);
 3.37-3.27 (1H, multiplet);
 3.00-2.90 (2H, multiplet);
 2.23-2.10 (2H, multiplet);
 1.96-1.83 (2H, multiplet);
 1.79-1.60 (2H, multiplet).

12(ii) (±)-α-(1-Benzylpiperidin-4-yl)amino-α-(pyridin-4-yl)acetonitrile

[0251] A mixture of 11.73 g (42.00 mmol) of 1-benzyl-4-(pyridin-4-yl)-methyleaminopiperidine [prepared as described in step 12(i) above] and 9.22 ml (65.67 mmol) of 95% trimethylsilyl cyanide was stirred at 100°C for 3 hours and then it was cooled to room temperature. 100 ml of methanol were added to the reaction mixture and the resulting mixture was then stirred for 30 minutes. At the end of this time, the solvent was distilled off from the reaction mixture by evaporation under reduced pressure to afford 12.87 g (quantitative yield) of the title compound as a brown oil.

[0252] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.67 (2H, doublet, J=6 Hz);
 7.50 (2H, doublet, J=6 Hz);
 7.37-7.22 (5H, multiplet);
 4.35 (1H, singlet);
 3.53 (2H, singlet);

2.98-2.77 (4H, multiplet);
2.33-1.44 (5H, multiplet).

12(iii) 1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0253] 14.25 g (46.51 mmol) of α -(1-benzylpiperidin-4-yl)amino- α -(pyridin-4-yl)acetonitrile [prepared as described in step 12(ii) above] and 6.98 g (46.51 mmol) of 3-(4-fluorophenyl)acrolein were dissolved in 145 ml of N,N-dimethylacetamide, 1.29 g (9.30 mmol) of potassium carbonate were added to the solution, and then this was stirred at room temperature for 5 hours. At the end of this time, water was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water and then concentrated by evaporation under reduced pressure. 150 ml of ethylene glycol were added to the residue thus obtained and the resulting mixture was stirred at 180°C for 1 hour. After cooling to room temperature, a saturated aqueous solution of sodium hydrogencarbonate was added to the reaction mixture which was then extracted with ethyl acetate. The organic extract was washed with water and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 1:3 by volume mixture of hexane and ethyl acetate as the eluant to give 10.31 g (yield 49%) of the title compound as a pale brown powder.

[0254] Melting point: 132 - 139°C

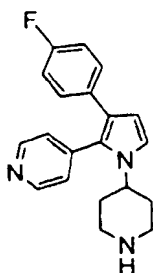
[0255] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.61 (2H, doublet, J=6 Hz);
7.37-7.23 (5H, multiplet);
7.13 (2H, doublet, J=6 Hz);
7.03 (2H, doublet of doublets, J=9 Hz, 5 Hz);
6.96 (1H, doublet, J=3 Hz);
6.37 (2H, triplet, J=9 Hz);
6.38 (1H, doublet, J=3 Hz);
3.37-3.60 (1H, multiplet);
3.50 (2H, singlet);
2.97 (2H, doublet, J=12 Hz);
2.12-1.56 (6H, multiplet).

Example 13

3-(4-Fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-142)

[0256]



[0257] In a similar manner to that described in Example 9 (ii) above, 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 12 above) was debenzylated to give the title compound (yield 30%) as a white powder.

[0258] Melting point: 191 - 192°C

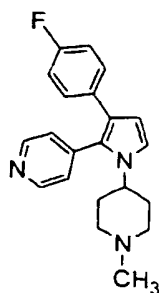
[0259] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.62 (2H, doublet, J=6 Hz);
7.15 (2H, doublet, J=6 Hz);
7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);
6.97 (1H, doublet, J=3 Hz);

- 6.38 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet, J=3 Hz);
 3.97-3.86 (1H, multiplet);
 3.17 (2H, doublet, J=12 Hz);
 2.58 (2H, doublet of triplets, J=12 Hz, 3 Hz);
 1.98-1.82 (4H, multiplet).

Example 14

- 3-(4-Fluorophenyl)-1-(1-methylpiperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-144)

[0260]

[0261] In a similar manner to the procedures described in Examples 4(i) and 4(ii) above, 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 13 above) was subjected consecutively to N-t-butoxycarbonylation and reduction to give the title compound (total yield for the two steps: 55%) as a white powder.

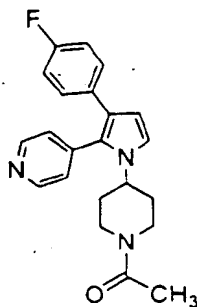
[0262] Melting point: 187- 189°C

[0263] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

- 8.62 (2H, doublet, J=6 Hz);
 7.14 (2H, doublet, J=6 Hz);
 7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.96 (1H, doublet, J=3 Hz);
 6.37 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet, J=3 Hz);
 3.37-3.77 (1H, multiplet);
 2.93 (2H, doublet, J=12 Hz);
 2.29 (3H, singlet);
 2.13-1.85 (6H, multiplet).

Example 151-(1-Acetylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2524)

[0264]



[0265] In a similar manner to that described in Example 3 above, a reaction was carried out using 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 13 above) to give the title compound (yield 98%) as a white powder.

[0266] Melting point: 223 - 224°C

[0267] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.64 (2H, doublet, J=6 Hz);

7.15 (2H, doublet, J=6 Hz);

7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.38 (2H, triplet, J=9 Hz);

6.37 (1H, doublet, J=3 Hz);

6.41 (1H, doublet, J=3 Hz);

4.34-4.74 (1H, multiplet);

4.05 (1H, doublet of triplets, J=12 Hz, 4 Hz);

3.95-3.83 (1H, multiplet);

3.08-2.97 (1H, multiplet);

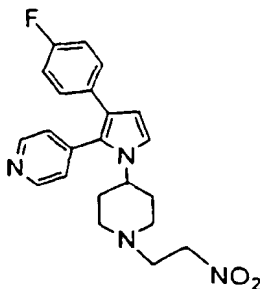
2.58-2.48 (1H, multiplet);

2.13 (3H, singlet);

2.04-1.78 (4H, multiplet).

Example 163-(4-Fluorophenyl)-1-[1-(2-nitroethyl)piperidin-4-yl]-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2537)

[0268]



[0269] 186 mg (1.40 mmol) of 2-nitroethyl acetate were added to a solution of 300 mg (0.93 mmol) of 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1*H*-pyrrole (prepared as described in Example 13 above) in 6 ml of ethanol. The resulting mixture was stirred at room temperature for 30 minutes, after which the solvent was distilled off under reduced pressure. A saturated aqueous solution of sodium hydrogencarbonate was added to the residue and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residual solid was washed with diethylether to give 280 mg (yield 76%) of the title compound as a white powder.

[0270] Melting point: 172- 173°C

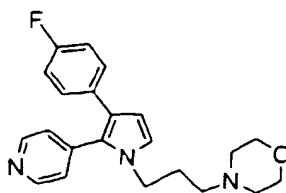
[0271] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.61 (2H, doublet, J=6 Hz);
 7.13 (2H, doublet, J=6 Hz);
 7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.93 (1H, doublet, J=3 Hz);
 6.37 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet, J=3 Hz);
 4.47 (2H, triplet, J=6 Hz);
 3.32 (1H, doublet of triplets, J=12 Hz, 4 Hz);
 3.03-2.92 (4H, multiplet);
 2.16-1.85 (6H, multiplet).

Example 17

3-(4-Fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1*H*-pyrrole (Compound No. 2-125)

[0272]



[0273] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 1-(3-aminopropyl)morpholine instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 13%) as pale brown needle-like crystals.

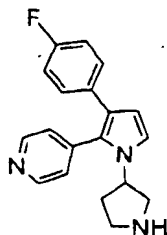
[0274] Melting point: 111 - 112°C

[0275] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.59 (2H, doublet, J=6 Hz);
 7.18 (2H, doublet, J=6 Hz);
 7.06 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.39 (2H, triplet, J=9 Hz);
 6.34 (1H, doublet, J=3 Hz);
 6.34 (1H, doublet, J=3 Hz);
 3.96 (2H, triplet, J=7 Hz);
 3.64 (2H, triplet, J=5 Hz);
 2.33-2.22 (4H, multiplet);
 2.20 (2H, triplet, J=7 Hz);
 1.72 (2H, quintet, J=7 Hz).

Example 18(±)-3-(4-Fluorophenyl)-2-(pyridin-4-yl)-1-(pyrrolidin-3-yl)-1H-pyrrole (Compound No. 2-139)

[0276]

18(i) (±)-1-[1-(t-Butoxycarbonyl)pyrrolidin-3-yl]-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0277] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using (±)-3-amino-1-(t-butoxycarbonyl)pyrrolidine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 31%) as a brown amorphous solid.

[0278] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.63 (2H, doublet, J=6 Hz);
 7.17 (2H, doublet, J=6 Hz);
 7.03 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.93-6.84 (3H, multiplet);
 6.41 (1H, doublet, J=3 Hz);
 4.70-4.55 (1H, multiplet);
 3.30-3.36 (4H, multiplet);
 2.30-2.12 (2H, multiplet);
 1.47 (9H, singlet).

18(ii) (±)-3-(4-Fluorophenyl)-2-(pyridin-4-yl)-1-(pyrrolidin-3-yl)-1H-pyrrole

[0279] 6.01 ml (24.05 mmol) of a 4N solution of hydrogen chloride in dioxane were added to a solution of 1.96 g (4.81 mmol) of (±)-1-[1-(t-butoxycarbonyl)pyrrolidin-3-yl]-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 18(i) above] in 10 ml of ethanol. The resulting mixture was stirred at 50°C for 30 minutes, after which the solvent was distilled off under reduced pressure. A saturated aqueous solution of sodium hydrogencarbonate was added to the residue thus obtained, and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure to give 1.29 g (yield 87%) of the title compound as a pale yellow powder.

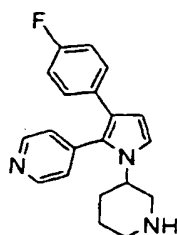
[0280] Melting point: 145 - 147°C

[0281] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.61 (2H, doublet, J=6 Hz);
 7.16 (2H, doublet, J=6 Hz);
 7.05 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.00 (1H, doublet, J=3 Hz);
 6.38 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet, J=3 Hz);
 4.61-4.53 (1H, multiplet);
 3.30-3.17 (2H, multiplet);
 3.10-2.97 (2H, multiplet);
 2.28-2.18 (2H, multiplet);
 2.11-2.01 (2H, multiplet).

Example 19(±)-3-(4-Fluorophenyl)-1-(piperidin-3-yl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-141)

[0282]



[0283] In a similar manner to the procedures described in Examples 18(i) and 18(ii) above, reactions were carried out using (±)-3-amino-1-(t-butoxycarbonyl)piperidine as a starting material instead of (±)-3-amino-1-(t-butoxycarbonyl)pyrrolidine to give the title compound (total yield 21%) as a white powder.

[0284] Melting point 189 - 191°C

[0285] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.61 (2H, doublet, J=6 Hz);

7.15 (2H, doublet, J=6 Hz);

7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.93 (1H, doublet, J=3 Hz);

6.37 (2H, triplet, J=9 Hz);

6.38 (1H, doublet, J=3 Hz);

3.93 (1H, triplet of triplets, J=11 Hz, 4 Hz);

3.20-3.12 (1H, multiplet);

3.05-2.96 (1H, multiplet);

2.33 (1H, triplet, J=11 Hz);

2.58 (1H, doublet of triplets, J=12 Hz, 3 Hz);

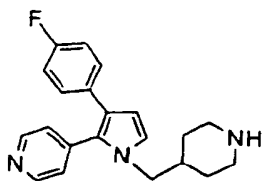
2.17-2.08 (1H, multiplet);

1.90-1.76 (2H, multiplet);

1.58-1.43 (1H, multiplet).

Example 203-(4-Fluorophenyl)-1-(piperidin-4-yl)methyl-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-169)

[0286]



[0287] In a similar manner to the procedures described in Examples 18(i) and 18(ii) above, reactions were carried out using 4-aminomethyl-1-(t-butoxycarbonyl)-piperidine as a starting material instead of 3-amino-1-(t-butoxycarbonyl)pyrrolidine to give the title compound (total yield 4%) as a white powder.

[0288] Melting point 174 - 176°C

[0289] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

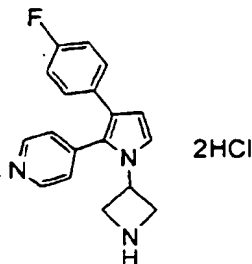
8.59 (2H, doublet, J=6 Hz);

7.15 (2H, doublet, J=6 Hz);

7.06 (2H, doublet of doublets J=9 Hz, 5 Hz);
 6.38 (2H, triplet, J=9 Hz);
 6.79 (1H, doublet, J=3 Hz);
 6.34 (1H, doublet, J=3 Hz);
 3.75 (2H, doublet, J=7 Hz);
 3.05-2.95 (2H, multiplet);
 2.45 (2H, doublet of triplets, J=12 Hz, 2 Hz);
 1.62-1.52 (1H, multiplet);
 1.46-1.38 (2H, multiplet);
 1.05-0.93 (2H, multiplet).

Example 21

1-(Azetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole dihydrochloride (Dihydrochloride of Compound No. 2-136)

[0290]

21(i) 1-(1-Diphenylmethylazetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0291] In a similar manner to the procedures in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 3-amino-1-diphenylmethylazetidine as a starting material instead of 4-amino-1-benzylpiperidine to afford the title compound (total yield 20%) as a pale brown powder.

[0292] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.56 (2H, doublet, J=6 Hz);
 7.43-7.18 (11H, multiplet);
 7.07-7.02 (4H, multiplet);
 6.38 (2H, triplet, J=9 Hz);
 6.43 (1H, doublet, J=3 Hz);
 4.74-4.66 (1H, multiplet);
 4.41 (1H, singlet);
 3.58-3.52 (2H, multiplet);
 3.34-3.28 (2H, multiplet).

21(ii) 1-(Azetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole dihydrochloride

[0293] 0.45 ml of concentrated hydrochloric acid and 0.12g of 20% palladium hydroxide on carbon were added to a solution of 1.20 g (2.61 mmol) of 1-(1-diphenylmethylazetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 21(i) above] in 12 ml of ethanol. The resulting mixture was stirred under a hydrogen atmosphere at 50°C for 20 hours. After cooling to room temperature, the reaction mixture was then filtered and the filtrate was concentrated by evaporation under reduced pressure. The residual solid was washed with diethylether to give 0.91 g (yield 95%) of the title compound as a pale yellow amorphous solid.

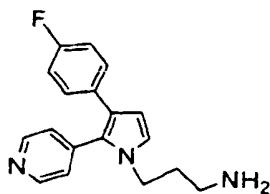
[0294] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

9.34-9.64 (1H, broad singlet);
 9.58-9.37 (1H, broad singlet);
 9.80 (2H, doublet, J=6 Hz);

7.94 (1H, doublet, J=3 Hz);
 7.64 (2H, doublet, J=6 Hz);
 7.25 (2H, triplet, J=9 Hz);
 7.20-7.07 (4H, multiplet);
 5 7.62 (1H, doublet, J=3 Hz);
 5.19-5.08 (1H, multiplet);
 4.43-4.23 (4H, multiplet).

Example 22

10 1-(3-Aminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-97)

[0295]

25 22(i) 1-(3-Benzyloxypropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0296] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii), reactions were carried out using 3-benzyloxypropylamine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 34%) as a brown oil.

30 **[0297]** ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.55 (2H, doublet, J=6 Hz);
 7.38-7.23 (5H, multiplet);
 7.16 (2H, doublet, J=6 Hz);
 7.07 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 35 6.89 (2H, triplet, J=9 Hz);
 6.31 (1H, doublet, J=3 Hz);
 6.34 (1H, doublet, J=3 Hz);
 4.41 (2H, singlet);
 4.03 (2H, triplet, J=7 Hz);
 40 3.36 (2H, triplet, J=7 Hz);
 1.86 (2H, quintet, J=7 Hz).

22(ii) 3-(4-Fluorophenyl)-1-(3-hydroxypropyl)-2-(pyridin-4-yl)-1H-pyrrole

45 **[0298]** In a similar manner to that described in Example 9(ii) above, 1-(3-benzyloxypropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 22(i) above] was debenzylated to afford the title compound (yield 55%) as a white powder.

[0299] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

50 8.59 (2H, doublet, J=6 Hz);
 7.19 (2H, doublet, J=6 Hz);
 7.05 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.89 (2H, triplet, J=9 Hz);
 6.37 (1H, doublet, J=3 Hz);
 6.36 (1H, doublet, J=3 Hz);
 55 4.05 (2H, triplet, J=7 Hz);
 3.55 (2H, triplet, J=7 Hz);
 1.30 (2H, quintet, J=7 Hz).

22(iii) 3-(4-Fluorophenyl)-2-(pyridin-4-yl)-1-[3-(p-toluenesulfonyloxy)propyl]-1H-pyrrole

[0300] 1.99 g (6.09 mmol) of p-toluenesulfonic anhydride and 0.95 ml (6.09 mmol) of triethylamine were added to a solution of 1.64 g (5.53 mmol) of 3-(4-fluorophenyl)-1-(3-hydroxypropyl)-2-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 22(ii) above] in 16 ml of dichloromethane. The resulting mixture was stirred at room temperature for 2 hours, after which the solvent was distilled off by evaporation under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using a 1:1 by volume mixture of hexane and ethyl acetate as the eluant to afford 1.33 g (yield 53%) of the title compound as a yellow powder.

[0301] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.57 (2H doublet, J=6 Hz);
 7.74 (2H doublet, J=8 Hz);
 7.35 (2H doublet, J=8 Hz);
 7.10 (2H doublet, J=6 Hz);
 7.03 (2H doublet of doublets, J=9 Hz, 5 Hz);
 6.88 (2H triplet, J=9 Hz);
 6.71 (1H doublet, J=3 Hz);
 6.29 (1H doublet, J=3 Hz);
 3.98 (2H triplet, J=7 Hz);
 3.39 (2H triplet, J=7 Hz);
 2.46 (3H singlet);
 1.34 (2H quintet, J=7 Hz).

22(iv) 1-(3-Azidopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0302] 0.38 g (5.90 mmol) of sodium azide were added to a solution of 1.33 g (2.95 mmol) of 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-[3-(p-toluenesulfonyloxy)propyl]-1H-pyrrole [prepared as described in step 22(iii) above] in 13 ml of dimethylsulfoxide. The resulting mixture was stirred at 70°C for 1 hour. At the end of this time water was added to the reaction mixture and then this was extracted with dichloromethane. The organic extract was washed with water and then concentrated under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 1:1 by volume mixture of hexane and ethylacetate as the eluant to afford 100 mg (yield 11%) of the title compound as a pale brown powder.

[0303] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.16 (2H doublet, J=6 Hz);
 7.17 (2H doublet, J=6 Hz);
 7.07 (2H doublet of doublets, J=9 Hz, 5 Hz);
 6.39 (2H triplet, J=9 Hz);
 6.33 (1H doublet, J=3 Hz);
 6.37 (1H doublet, J=3 Hz);
 4.00 (2H triplet, J=7 Hz);
 3.18 (2H triplet, J=7 Hz);
 1.78 (2H quintet, J=7 Hz).

22(v) 1-(3-Aminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole

[0304] 20 mg of 10% palladium on carbon were added to a solution of 100 mg (0.31 mmol) of 1-(3-azidopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole [prepared as described in Step 22(iv) above] in 2 ml of ethanol. The resulting mixture was stirred under a hydrogen atmosphere at room temperature for 1 hour, after which the reaction mixture was filtered and the filtrate was concentrated by evaporation under reduced pressure. The residual solid was washed with diethylether to give 43 mg (yield 47%) of the title compound as a pale yellow powder.

[0305] Melting point: 219 - 222°C

[0306] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

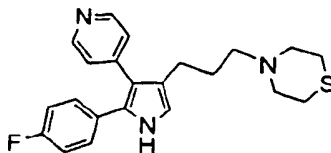
8.61 (2H doublet, J=6 Hz);
 7.74-7.40 (2H, broad singlet);
 7.27 (2H doublet, J=6 Hz);
 7.12-7.00 (5H, multiplet);
 6.38 (1H doublet, J=3 Hz);
 3.97 (2H triplet, J=7 Hz);
 2.60 (2H triplet, J=7 Hz);

1.73 (2H, quintet, J=7 Hz).

Example 23

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-[3-(thiomorpholin-1-yl)propyl]-1H-pyrrole (Compound No. 1-129)

[0307]



23(i) Ethyl 3-[2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrol-4-yl]acrylate

[0308] In a similar manner to that described in Example 1(iv) above, a reaction was carried out using ethyl diethylphosphonoacetate instead of diethylphosphonoacetonitrile to afford the title compound (yield 60%) as a brown powder.

[0309] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.64 (1H, broad singlet);
8.56 (2H, doublet, J=6 Hz);
7.54 (1H, doublet, J=16 Hz);
7.16 (2H, doublet, J=6 Hz);
7.16-7.13 (3H, multiplet);
6.99 (2H, triplet, J=9 Hz);
6.07 (1H, doublet, J=16 Hz);
4.20 (2H, quartet, J=7 Hz);
1.28 (3H, triplet, J=7 Hz).

23(ii) 4-(2-Ethoxycarbonyl)ethyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0310] In a similar manner to that described in Example 1(v) above, ethyl 3-[2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrol-4-yl]acrylate [prepared as described in step 23(i) above] was reduced to afford the title compound (yield 73%) as a pale brown powder.

[0311] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.51 (2H, doublet, J=4 Hz);
8.33 (1H, broad singlet);
7.15 (2H, doublet, J=4 Hz);
7.16-7.12 (2H, multiplet);
6.96 (2H, triplet, J=9 Hz);
6.74 (1H, doublet, J=3 Hz);
4.11 (2H, quartet, J=7 Hz);
2.85 (2H, triplet, J=6 Hz);
2.50 (2H, triplet, J=6 Hz);
1.23 (3H, triplet, J=7 Hz).

23(iii) 2-(4-Fluorophenyl)-4-(3-hydroxypropyl)-3-(pyridin-4-yl)-1H-pyrrole

[0312] In a similar manner to that described in Example 1(vi) above, 4-(2-ethoxycarbonyl)ethyl-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 23(ii) above] was reduced using lithium aluminum hydride to afford the title compound (yield 91%) as a pale brown powder.

[0313] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.51 (2H, doublet, J=5 Hz);
8.23 (1H, broad singlet);
7.17-7.12 (4H, multiplet);
7.00 (2H, triplet, J=9 Hz);

6.75 (1H, doublet, J=3 Hz);
 3.65 (2H, triplet, J=6 Hz);
 2.61 (2H, triplet, J=6 Hz);
 1.79-1.72 (2H, multiplet);
 1.61 (1H, broad singlet).

23(iv) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-[3-(p-toluenesulfonyloxy)propyl]-1H-pyrrole

[0314] In a similar manner to that described in Example 22(iii) above, p-toluenesulfonylation was carried out using 2-(4-fluorophenyl)-4-(3-hydroxypropyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 23(iii) above] to afford the title compound (yield 71%) as a pale yellow powder.

[0315] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.49 (2H, doublet, J=6 Hz);
 8.20 (1H, broad singlet);
 7.59 (2H, doublet, J=8 Hz);
 7.33 (2H, doublet, J=8 Hz);
 7.12 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.08 (2H, doublet, J=6 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.69 (1H, doublet, J=3 Hz);
 4.00 (2H, triplet, J=6 Hz);
 2.56 (2H, triplet, J=6 Hz);
 2.44 (3H, singlet);
 1.31-1.74 (2H, multiplet).

23(v) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-[3-(thiomorpholin-1-yl)propyl]-1H-pyrrole

[0316] 163 mg (1.18 mmol) of potassium carbonate and 122 μl (1.29 mmol) of thiomorpholine were added to a solution of 450 mg (1.07 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-[3-(p-toluenesulfonyloxy)propyl]-1H-pyrrole [prepared as described in step 23(iv) above] in 50 ml of acetonitrile. The resulting mixture was heated under reflux overnight. After cooling the reaction mixture to room temperature, water was added and then this was extracted with ethyl acetate. The organic extract was washed with water and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 95:5 by volume mixture of ethyl acetate and methanol as the eluant to give 340 mg (yield 83%) of the title compound as a pale yellow powder.

[0317] Melting point: 205 - 207°C

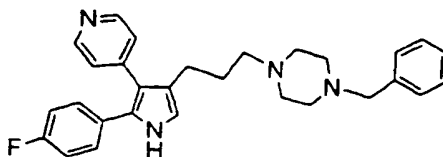
[0318] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.51 (2H, doublet, J=6 Hz);
 8.20 (1H, broad singlet);
 7.14 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.13 (2H, doublet, J=6 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.72 (1H, doublet, J=3 Hz);
 2.65 (8H, singlet);
 2.50 (2H, triplet, J=6 Hz);
 2.35 (2H, triplet, J=6 Hz);
 1.66 (2H, quintet, J=8 Hz).

Example 24

4-[3-(1-Benzylpiperazin-4-yl)propyl]-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-135)

[0319]



[0320] In a similar manner to that described in Example 23(v) above, a reaction was carried out using 1-benzylpiperazine instead of thiomorpholine to give the title compound (yield 52%) as a white powder.

[0321] Melting point: 153 - 154°C

[0322] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.50 (2H, doublet, J=5 Hz);

8.28 (1H, broad singlet);

7.31 (4H, doublet, J=4 Hz);

7.26-7.22 (1H, multiplet);

7.13 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.13 (2H, doublet, J=5 Hz);

6.96 (2H, triplet, J=9 Hz);

6.72 (1H, doublet, J=2 Hz);

3.50 (2H, singlet);

2.50 (2H, triplet, J=8 Hz);

2.44 (8H, broad singlet);

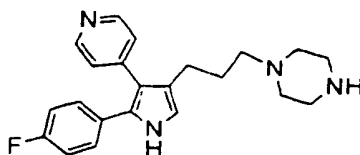
2.34 (2H, triplet, J=8 Hz);

1.69 (2H, quintet, J=8 Hz).

Example 25

2-(4-Fluorophenyl)-4-[3-(piperazin-1-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-133)

[0323]



[0324] In a similar manner to that described in Example 9(ii) above, 4-[3-(1-benzylpiperazin-4-yl)propyl]-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 24 above) was debenzylated to give the title compound (yield 98%) as a white powder.

[0325] Melting point: 162 - 164°C

[0326] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.51 (2H, doublet, J=6 Hz);

8.20 (1H, broad singlet);

7.14 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.14 (2H, doublet, J=6 Hz);

6.96 (2H, triplet, J=9 Hz);

6.73 (1H, doublet, J=2 Hz);

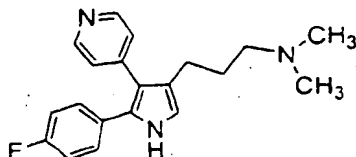
2.37 (4H, triplet, J=5 Hz);

2.51 (2H, triplet, $J=8$ Hz);
 2.36 (4H, broad singlet);
 2.32 (2H, triplet, $J=8$ Hz);
 1.70 (2H, quintet, $J=8$ Hz).

Example 26

4-(3-Dimethylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-109)

[0327]



[0328] In a similar manner to that described in Example 23(v) above, a reaction was carried out using dimethylamine instead of thiomorpholine to give the title compound (yield 66%) as a white powder.

[0329] Melting point: 177 - 179°C

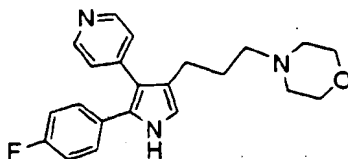
[0330] ^1H -Nuclear magnetic resonance spectrum (500 MHz, CDCl_3) δ ppm:

8.51 (2H, doublet, $J=5$ Hz);
 8.24 (1H, broad singlet);
 7.15-7.13 (2H, multiplet);
 7.14 (2H, doublet, $J=5$ Hz);
 6.96 (2H, triplet, $J=9$ Hz);
 6.74 (1H, doublet, $J=2$ Hz);
 2.52 (2H, triplet, $J=8$ Hz);
 2.28 (2H, triplet, $J=8$ Hz);
 2.19 (6H, singlet);
 1.74-1.59 (2H, multiplet).

Example 27

2-(4-Fluorophenyl)-4-[3-(morpholin-1-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-125)

[0331]



[0332] In a similar manner to that described in Example 23(v) above, a reaction was carried out using morpholine instead of thiomorpholine to give the title compound (yield 45%) as a pale yellow powder.

[0333] Melting point: 182 - 183°C

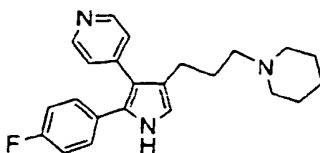
[0334] ^1H -Nuclear magnetic resonance spectrum (400 MHz, CDCl_3) δ ppm:

8.62 (1H, broad singlet);
 8.50 (2H, doublet, $J=5$ Hz);
 7.17-7.12 (2H, multiplet);
 7.14 (2H, doublet, $J=5$ Hz);
 6.96 (2H, triplet, $J=9$ Hz);

6.73 (1H, doublet, J=3 Hz);
 3.70-3.68 (4H, multiplet);
 2.53 (2H, triplet, J=8 Hz);
 2.41-2.34 (4H, multiplet);
 2.33 (2H, triplet, J=8 Hz);
 1.68 (2H, quintet, J=8 Hz).

Example 28

2-(4-Fluorophenyl)-4-[3-(piperidin-1-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-121)

[0335]

[0336] In a similar manner to that described in Example 23(v) above, a reaction was carried out using piperidine instead of thiomorpholine to give the title compound (yield 63%) as a pale yellow powder.

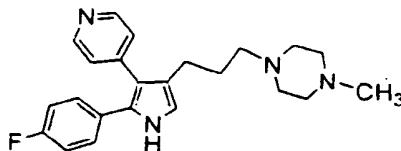
[0337] Melting point: 195 - 197°C

[0338] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.50 (2H, doublet, J=6 Hz);
 8.19 (1H, broad singlet);
 7.14 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.74 (1H, doublet, J=2 Hz);
 2.50 (2H, triplet, J=8 Hz);
 2.40-2.30 (4H, multiplet);
 1.75-1.55 (8H, multiplet);
 1.43-1.42 (2H, multiplet).

Example 29

2-(4-Fluorophenyl)-4-[3-(1-methylpiperazin-4-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-134)

[0339]

[0340] In a similar manner to that described in Example 23(v) above, a reaction was carried out using 1-methylpiperazine instead of thiomorpholine to give the title compound (yield 64%) as a pale yellow powder.

[0341] Melting point: 190 - 192°C

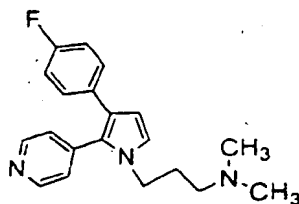
[0342] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.65 (1H, broad singlet);
 8.49 (2H, doublet, J=6 Hz);
 7.16-7.12 (2H, multiplet);
 7.14 (2H, doublet, J=6 Hz);
 6.95 (2H, triplet, J=9 Hz);

6.73 (1H, doublet, J=3 Hz);
 2.52 (2H, triplet, J=8 Hz);
 2.53-2.36 (8H, multiplet);
 2.36 (2H, triplet, J=8 Hz);
 2.28 (3H, singlet);
 1.70 (2H, quintet, J=8 Hz).

Example 30

1-(3-Dimethylaminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-109)

[0343]

[0344] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 3-dimethylaminopropylamine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 9%) as a white powder.

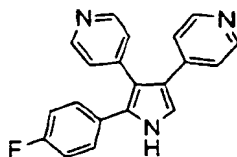
[0345] Melting point: 93 - 95°C

[0346] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.59 (2H, doublet, J=6 Hz);
 7.18 (2H, doublet, J=6 Hz);
 7.07 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 6.39 (2H, triplet, J=9 Hz);
 6.34 (1H, doublet, J=3 Hz);
 6.35 (1H, doublet, J=3 Hz);
 3.93 (2H, triplet, J=7 Hz);
 2.14 (2H, triplet, J=7 Hz);
 2.11 (6H, singlet);
 1.71 (2H, quintet, J=7 Hz).

Example 31

2-(4-Fluorophenyl)-3,4-bis(pyridin-4-yl)-1H-pyrrole (Compound No. 1-159)

[0347]

[0348] 450 mg (1.15 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole dihydrochloride [prepared as described in Example 10 above] were suspended in a mixture of 25 ml of xylene and 2 ml of methanol and 450 mg of 10% palladium on carbon and 320 mg (2.30 mmol) of sodium carbonate were added to the suspension. The resulting mixture was heated under reflux for 4 hours. At the end of this time, the reaction mixture was filtered and the resulting filtrate was concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 10:1:1 by volume mixture of ethyl acetate, meth-

anol and isopropylamine as the eluant to give 192 mg (yield 53%) of the title compound as a pale yellow powder.

[0349] Melting point: 325 - 330°C (decomposition)

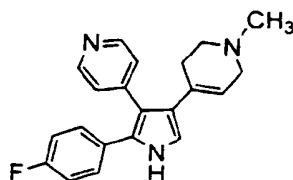
[0350] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

- 11.87 (1H, broad singlet);
 8.47 (2H, doublet, J=6 Hz);
 8.35 (2H, doublet, J=6 Hz);
 7.39 (1H, doublet, J=3 Hz);
 7.21 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.16 (2H, triplet, J=9 Hz);
 7.08 (2H, doublet, J=6 Hz);
 7.03 (2H, doublet, J=6 Hz).

Example 32

15 2-(4-Fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-145)

[0351]



30 32(i) 2-(4-Fluorophenyl)-4-(4-hydroxy-1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0352] In a similar manner to that described in Example 7(iv) above, a reaction was carried out using 1-methylpiperidine-4-one instead of pentafluoropyridine to afford the title compound (yield 47%) as a pale brown powder.

[0353] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

- 8.22 (2H, doublet, J=6 Hz);
 7.27 (2H, doublet, J=6 Hz);
 7.21 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.94 (2H, triplet, J=9 Hz);
 6.92 (1H, singlet);
 2.56-2.42 (4H, broad multiplet);
 2.22 (3H, singlet);
 2.00-1.91 (2H, broad multiplet);
 1.34-1.78 (2H, broad multiplet);
 1.20-1.11 (3H, multiplet);
 1.06 (16H, doublet, J=7 Hz).

45 32(ii) 2-(4-Fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole

[0354] In a similar manner to that described in Example 9(iii) above, using 2-(4-fluorophenyl)-4-(4-hydroxy-1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 32(i) above], dehydration with triethylsilane/trifluoroacetic acid and desilylation with tetrabutylammonium fluoride were carried out to give the title compound (yield 91%) as a pale yellow powder.

[0355] Melting point: 230 - 232°C (decomposition)

[0356] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

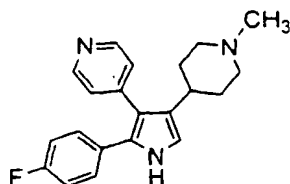
- 11.36 (1H, broad singlet);
 8.45 (2H, doublet, J=6 Hz);
 7.19-7.14 (6H, multiplet);
 6.90 (1H, doublet, J=3 Hz);
 5.24-5.21 (1H, broad multiplet);

2.78-2.75 (2H, broad multiplet);
2.40 (2H, triplet, J=5 Hz);
2.18 (3H, singlet);
2.20-2.13 (2H, broad multiplet).

Example 33

2-(4-Fluorophenyl)-4-(1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-144)

[0357]



[0358] In a similar manner to that described in Example 11 above, 2-(4-fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 32 above] was reduced to give the title compound (yield 85%) as a white powder.

[0359] Melting point: 284 - 285°C (decomposition)

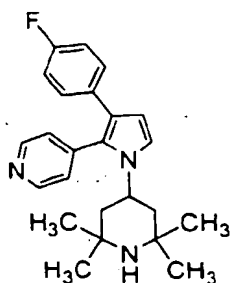
[0360] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.20 (1H, broad singlet);
8.48 (2H, doublet, J=6 Hz);
7.19-7.16 (4H, multiplet);
7.10 (2H, triplet, J=9 Hz);
6.76 (1H, doublet, J=2 Hz);
2.73 (2H, doublet, J=12 Hz);
2.35 (1H, triplet of triplets, J=12 Hz, 4 Hz);
2.12 (3H, singlet);
1.77 (2H, triplet, J=12 Hz);
1.61 (2H, doublet, J=12 Hz);
1.51 (2H, double doublet of triplets, J=13 Hz, 12 Hz, 4 Hz).

Example 34

3-(4-Fluorophenyl)-2-(pyridin-4-yl)-1-(2,2,6,6-tetramethylpiperidin-4-yl)-1H-pyrrole (Compound No. 2-2523)

[0361]



[0362] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 4-amino-2,2,6,6-tetramethylpiperidine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 12%) as a white powder.

[0363] Melting point: 118 - 119°C

[0364] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.62 (2H, doublet, J=6 Hz);

7.18 (2H, doublet, J=6 Hz);

5 7.07 (2H, doublet of doublets, J=9 Hz, 6 Hz);

6.39 (1H, doublet, J=3 Hz);

6.38 (2H, triplet, J=9 Hz);

6.39 (1H, doublet, J=3 Hz);

4.43-4.38 (1H, multiplet);

10 1.36 (2H, doublet of doublets, J=13 Hz, 3 Hz);

1.55-1.38 (3H, multiplet);

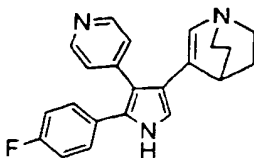
1.15 (6H, singlet);

1.05 (6H, singlet).

15 Example 35

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1H-pyrrole (Compound No. 1-2540)

[0365]



30 35(i) (+)-2-(4-Fluorophenyl)-4-(3-hydroxyquinuclidin-3-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0366] In a similar manner to that described in Example 7(iv) above, a reaction was carried out using quinuclidine-3-one instead of pentafluoropyridine to afford the title compound (yield 37%) as a pale brown powder.

[0367] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

35 8.25 (2H, doublet, J=6 Hz);

7.25 (2H, doublet, J=6 Hz);

7.26-7.16 (2H, broad multiplet);

7.08 (2H, triplet, J=9 Hz);

6.30 (1H, singlet);

40 4.32 (1H, singlet);

2.31 (1H, doublet, J=14 Hz);

2.72-2.64 (1H, broad multiplet);

2.59 (1H, doublet, J=15 Hz);

2.60-2.50 (2H, broad multiplet);

45 2.46-2.37 (1H, broad multiplet);

2.03-1.91 (2H, broad multiplet);

1.49-1.37 (2H, broad multiplet);

1.26-1.17 (1H, broad multiplet);

1.15-1.05 (3H, multiplet);

50 1.03-0.95 (18H, multiplet).

35(ii) 2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1H-pyrrole

[0368] 24 ml of formic acid were added to 1.20 g (2.31 mmol) of (±)-2-(4-fluorophenyl)-4-(3-hydroxyquinuclidin-3-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 35(i) above]. The resulting mixture was stirred at 90°C for 3 hours. After cooling to room temperature, water was added to the reaction mixture. The pH of the mixture was adjusted to greater than 10 through the addition of a 10% aqueous solution of sodium hydroxide and then the resulting mixture was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous

sodium sulfate and then concentrated by evaporation under reduced pressure. The residual solid was washed with a 1:1 by volume mixture of ethyl acetate and diethyl ether to give 568 mg (yield 71%) of the title compound as a pale brown powder.

[0369] Melting point: 246 - 248 (decomposition)

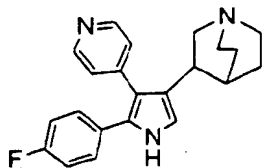
[0370] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.48 (1H, broad singlet);
8.47 (2H, doublet, J=6 Hz);
7.19-7.14 (4H, multiplet);
7.11 (2H, triplet, J=9 Hz);
7.00 (1H, doublet, J=2 Hz);
5.86 (1H, singlet);
2.77 (2H, double doublet of doublets, J=13 Hz, 9 Hz, 5 Hz);
2.61-2.56 (1H, broad multiplet);
2.40-2.31 (2H, multiplet);
1.58-1.49 (2H, broad multiplet);
1.43-1.34 (2H, broad multiplet).

Example 36

(±)-2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-3-yl)-1H-pyrrole (Compound No. 1-2539)

[0371]



[0372] In a similar manner to that described in Example 11 above, 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1H-pyrrole (prepared as described in Example 35 above) was reduced to give the title compound (yield 82%) as a white powder

[0373] Melting point: 239 - 241°C (decomposition)

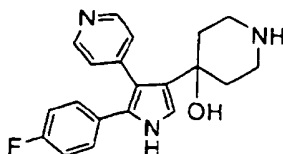
[0374] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.25 (1H, broad singlet);
8.47 (2H, doublet, J=6 Hz);
7.18-7.07 (6H, multiplet);
6.93 (1H, doublet, J=2 Hz);
2.98-2.90 (1H, multiplet);
2.89-2.76 (2H, broad multiplet);
2.74-2.58 (4H, broad multiplet);
1.73-1.64 (1H, broad multiplet);
1.54-1.46 (2H, broad multiplet);
1.42-1.33 (1H, broad multiplet);
1.26-1.17 (1H, broad multiplet).

Example 37

2-(4-Fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2526)

[0375]



[0376] 2.73 ml (2.73 mmol) of a 1M solution of tetrabutylammonium fluoride in tetrahydrofuran were added to a solution of 675 mg (1.37 mmol) of 2-(4-fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in Example 9(iii) above] in 14 ml of tetrahydrofuran and the resulting mixture was stirred at room temperature for 30 minutes. At the end of this time, the solvent was distilled off under reduced pressure. Water was added to the resulting residue and the precipitate thus obtained was collected by filtration and then washed with ethyl acetate to give 363 mg (yield 79%) of the title compound as a white powder.

[0377] Melting point: 197 - 199°C

[0378] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.16 (1H, broad singlet);

8.42 (2H, doublet, J=6 Hz);

7.32 (2H, doublet, J=6 Hz);

7.04-7.01 (4H, multiplet);

6.79 (1H, doublet, J=2 Hz);

4.45 (1H, singlet);

2.78-2.70 (2H, multiplet);

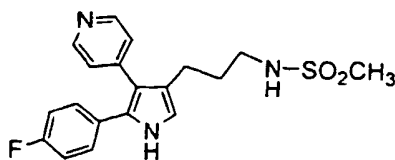
2.55-2.45 (2H, these overlapped with protons derived from remaining DMSO-d₆);

1.63-1.48 (4H, multiplet).

Example 38

2-(4-Fluorophenyl)-4-(3-methanesulfonylaminopropyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2536)

[0379]



[0380] 188 mg (1.08 mmol) of methanesulfonic anhydride were added to a solution of 290 mg (0.982 mmol) of 4-(3-aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 1 above) in 5 ml of pyridine and the resulting mixture was allowed to stand overnight. At the end of this time, water was added to the reaction mixture, giving a precipitate which was collected by filtration and washed with ethyl acetate to give 208 mg (yield 57%) of the title compound as a pale brown powder.

[0381] Melting point: 210 - 214°C

[0382] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.52 (2H, doublet, J=6 Hz);

8.23 (1H, broad singlet);

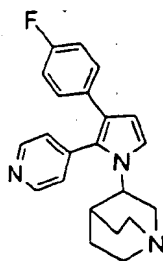
7.15-7.12 (4H, multiplet);

6.97 (2H, triplet, J=9 Hz);
 6.76 (1H, doublet, J=2 Hz);
 4.15 (1H, broad singlet);
 3.13-3.09 (2H, multiplet);
 2.89 (3H, singlet);
 2.61 (2H, triplet, J=7 Hz);
 1.72 (2H, quintet, J=7 Hz).

Example 39

(±)-3-(4-Fluorophenyl)-2-(pyridin-4-yl)-1-(quinuclidin-3-yl)-1H-pyrrole (Compound No. 2-2539)

[0383]



[0384] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using (±)-3-aminoquinuclidine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (total yield 25%) as a pale brown powder.

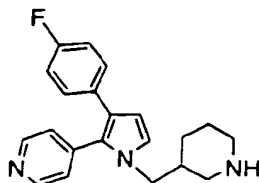
[0385] Melting point 229 - 230°C

[0386] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm: .

8.61 (2H, doublet, J=6 Hz);
 7.16 (2H, doublet, J=6 Hz);
 7.13 (1H, doublet, J=3 Hz);
 7.04 (2H, doublet of doublets J=9 Hz, 5 Hz);
 6.57 (2H, triplet, J=9 Hz);
 6.42 (1H, doublet, J=3 Hz);
 4.24-4.15 (1H, multiplet);
 3.33-3.23 (1H, multiplet);
 3.22-3.05 (2H, multiplet);
 2.98-2.79 (2H, multiplet);
 2.78-2.67 (1H, multiplet);
 1.95-1.81 (2H, multiplet);
 1.71-1.58 (1H, multiplet);
 1.56-1.37 (2H, multiplet).

Example 40(±)-3-(4-Fluorophenyl)-1-(piperidin-3-yl)methyl-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2520)

[0387]



[0388] In a similar manner to the procedures described in Examples 18(i) and 18(ii) above, reactions were carried out using (±)-3-aminomethyl-(1-t-butoxycarbonyl)piperidine as a starting material instead of (±)-3-amino-1-(t-butoxycarbonyl)pyrrolidine to give the title compound (total yield 7%) as a white powder.

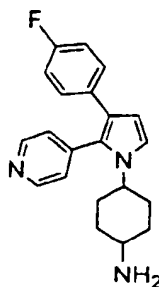
[0389] Melting point 134 - 135°C

[0390] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.59 (2H, doublet, J=6 Hz);
 7.16 (2H, doublet, J=6 Hz);
 7.06 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.38 (2H, triplet, J=9 Hz);
 6.79 (1H, doublet, J=3 Hz);
 6.34 (1H, doublet, J=3 Hz);
 3.74 (2H, doublet, J=8 Hz);
 2.94-2.87 (1H, multiplet);
 2.33-2.73 (1H, multiplet);
 2.46 (1H, doublet of doublets, J=12 Hz, 3 Hz);
 2.13 (1H, doublet of doublets, J=12 Hz, 10 Hz);
 1.75-1.51 (3H, multiplet);
 1.40-1.28 (1H, multiplet);
 0.98-0.85 (1H, multiplet).

Example 41cis-1-(4-Aminocyclohexyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2529)

[0391]



[0392] In a similar manner to the procedures described in Examples 18(i) and 18(ii) above, reactions were carried out using cis-4-(t-butoxycarbonylamino)cyclohexylamine as a starting material instead of (±)-3-amino-1-(t-butoxycarbonyl)pyrrolidine to give the title compound (total yield 10%) as a pale brown powder.

[0393] Melting point 163 - 164°C

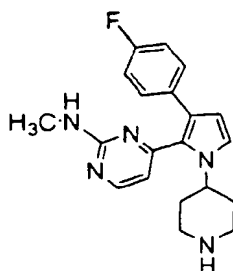
[0394] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.61 (2H, doublet, J=6 Hz);
 7.14 (2H, doublet, J=6 Hz);
 7.05 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 5 7.03 (1H, doublet, J=3 Hz);
 6.37 (2H, triplet, J=9 Hz);
 6.33 (1H, doublet, J=3 Hz);
 3.32 (1H, triplet of triplets, J=12 Hz, 4 Hz);
 3.29-3.22 (1H, multiplet);
 10 2.22-2.07 (2H, multiplet);
 1.32-1.41 (6H, multiplet).

Example 42

15 3-(4-Fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1-(piperidin-4-yl)-1H-pyrrole (Compound No. 2-1072)

[0395]



42(i) 1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylthiopyrimidin-4-yl)-1H-pyrrole

[0396] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 4-formyl-2-methylthiopyrimidine as a starting material instead of 4-formylpyridine to give the title compound (total yield 34%) as a brown oil.

[0397] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.21 (1H, doublet, J=5 Hz);
 7.39-7.24 (5H, multiplet);
 40 7.16 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.06 (1H, doublet, J=3 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.57 (1H, doublet, J=5 Hz);
 6.27 (1H, doublet, J=3 Hz);
 45 4.33-4.73 (1H, multiplet);
 3.54 (2H, singlet);
 3.02 (2H, doublet, J=9 Hz);
 2.75 (2H, triplet, J=6 Hz);
 2.46 (2H, triplet, J=6 Hz);
 50 2.13-1.96 (2H, multiplet).

42(ii) 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methanesulfonylpyrimidin-4-yl)-1H-pyrrole

[0398] 3.51 g (15.26 mmol) of 69-75 wt.% of m-chloroperbenzoic acid were added in small portions to a solution of 3.50 g (7.63 mmol) of 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylthiopyrimidin-4-yl)-1H-pyrrole [prepared as described in step 42(i) above] in 35 ml of ethyl acetate with ice-cooling. The resulting mixture was stirred at room temperature 2 days, after which a 10% aqueous solution of sodium thiosulfate was added to the reaction mixture followed by extraction with ethyl acetate. The organic extract was washed with water and then concentrated by evap-

oration under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 9:1 by volume mixture of dichloromethane and methanol as the eluant to give 2.48 g (yield 66%) of the title compound as a pale brown amorphous solid.

[0399] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.35 (1H, doublet, J=5 Hz);
 7.63-7.56 (2H, multiplet);
 7.47-7.38 (3H, multiplet);
 7.29-7.24 (1H, multiplet);
 7.22 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.05 (2H, triplet, J=9 Hz);
 6.90 (2H, doublet, J=5 Hz);
 6.29 (1H, doublet, J=3 Hz);
 5.43-5.33 (1H, multiplet);
 4.46 (2H, singlet);
 3.56-3.27 (4H, multiplet);
 3.03-2.86 (2H, multiplet);
 2.93 (3H, singlet);
 2.26-2.14 (1H, multiplet);
 2.10-2.01 (1H, multiplet).

42(iii) 1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1H-pyrrole

[0400] A mixture of 1.19 g (2.43 mmol) of 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methanesulfonylpyrimidin-4-yl)-1H-pyrrole [prepared as described in Step 42(ii) above] and 13 ml of a 2M solution of methylamine in tetrahydrofuran was stirred at 100°C for 30 minutes in a sealed tube. At the end of this time, the solvent was distilled off under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using a 1:2 by volume mixture of hexane and ethyl acetate as the eluant to give 260 mg (yield 24%) of the title compound as a yellow amorphous solid.

[0401] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.03 (1H, doublet, J=5 Hz);
 7.42-7.23 (5H, multiplet);
 7.18 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.00 (1H, doublet, J=3 Hz);
 6.94 (2H, triplet, J=9 Hz);
 6.26 (1H, doublet, J=3 Hz);
 6.23 (1H, doublet, J=5 Hz);
 5.14-5.04 (1H, multiplet);
 4.32-4.68 (1H, multiplet);
 3.53 (2H, singlet);
 3.10-2.95 (5H, multiplet);
 2.12-1.96 (6H, multiplet).

42(iv) 3-(4-Fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1-(piperidin-4-yl)-1H-pyrrole

[0402] In a similar manner to that described in Example 9(ii) above, 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1H-pyrrole [prepared as described in step 42(iii) above] was debenzylated to give the title compound (yield 39%) as a pale yellow powder.

[0403] Melting point: 189- 191°C

[0404] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

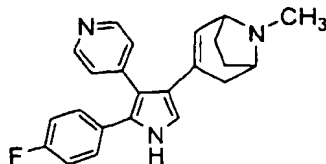
8.05 (1H, doublet, J=5 Hz);
 7.19 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.00 (1H, doublet, J=3 Hz);
 6.95 (2H, triplet, J=9 Hz);
 6.28 (1H, doublet, J=3 Hz);
 6.24 (1H, doublet, J=5 Hz);
 5.14-5.04 (1H, multiplet);
 4.96-4.81 (1H, multiplet);
 3.24 (2H, doublet, J=12 Hz);

3.04 (3H, doublet, J=5 Hz);
 2.70 (2H, doublet of quartets, J=12 Hz, 2 Hz);
 2.14 (2H, doublet, J=12 Hz);
 1.91 (2H, doublet of quartets, J=12 Hz, 4 Hz).

Example 43

(±)-2-(4-Fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2544)

[0405]



43(i) 2-(4-Fluorophenyl)-4-(3-hydroxy-8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0406] In a similar manner to that described in Example 7(iv) above, a reaction was carried out using 8-methyl-8-azabicyclo[3.2.1]octane-3-one instead of pentafluoropyridine to give a mixture of isomers of the title compound (yield 49%) as a pale brown powder.

[0407] From the mixture thus obtained, isomer A (Rf value = 0.45, pale brown powder) and isomer B (Rf value = 0.40, pale brown powder) were separated by chromatography on an alumina column using a 10:1 by volume mixture of ethyl acetate and methanol as the eluant (the ratio of isomer A:isomer B was 85:90 by weight).

isomer A

[0408] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.21 (2H, doublet, J=6 Hz);
 7.26 (2H, doublet, J=6 Hz);
 7.17 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.92 (2H, triplet, J=9 Hz);
 6.34 (1H, singlet);
 3.14-3.06 (1H, broad multiplet);
 2.28-2.15 (4H, multiplet);
 2.20 (3H, singlet);
 1.95-1.84 (4H, broad multiplet);
 1.19-1.09 (3H, multiplet);
 1.05 (1H, doublet, J=7 Hz).

isomer B

[0409] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.22 (2H, doublet, J=6 Hz);
 7.27 (2H, doublet, J=6 Hz);
 7.18 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.93 (2H, triplet, J=9 Hz);
 6.35 (1H, singlet);
 3.38 (1H, broad singlet);
 2.41 (3H, singlet);
 2.38-2.27 (4H, multiplet);
 2.06-1.95 (4H, broad multiplet);
 1.20-1.10 (3H, multiplet);

1.06 (1H, doublet, J=7 Hz).

43(ii) (±)-2-(4-Fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1H-pyrrole

[0410] In a similar manner to that described in Example 9(iii) above, 2-(4-fluorophenyl)-4-(3-hydroxy-8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1H-pyrrole (a mixture of isomers A and B, prepared as described in step 43(i) above) was subjected to dehydration using triethylsilane and trifluoroacetic acid followed by desilylation (deprotection) with tetrabutylammonium fluoride to give the title compound (yield 83%) as a white powder.

[0411] Melting point: 243 - 245°C

[0412] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.39 (2H, doublet, J=7 Hz);

7.25 (2H, doublet, J=6 Hz);

7.16 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.99 (2H, triplet, J=9 Hz);

6.33 (1H, singlet);

5.45 (1H, broad doublet, J=6 Hz);

3.28-3.24 (1H, broad multiplet);

3.22 (1H, triplet, J=6 Hz);

2.68 (1H, broad doublet, J=16 Hz);

2.37 (3H, singlet);

2.20-2.11 (1H, broad multiplet);

2.09-1.99 (1H, multiplet);

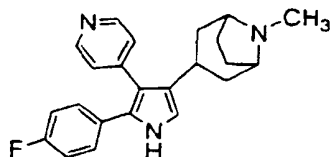
1.90-1.80 (2H, broad multiplet);

1.69-1.61 (1H, multiplet).

Example 44

2-(4-Fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2542)

[0413]



[0414] In a similar manner to that described in Example 11 above, (±)-2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 43 above] was reduced to give the title compound (yield 57%) as a white powder.

[0415] Melting point: 233 - 234°C (decomposition)

[0416] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.13 (1H, broad singlet);

8.49 (2H, doublet, J=6 Hz);

7.14-7.05 (6H, multiplet);

6.76 (1H, doublet, J=3 Hz);

3.03-2.65 (3H, broad multiplet);

2.16-2.07 (2H, multiplet);

2.04 (3H, singlet);

1.98-1.90 (2H, multiplet);

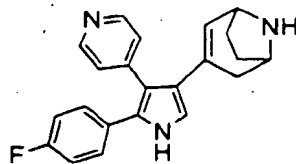
1.45-1.36 (2H, multiplet);

1.24-1.14 (2H, multiplet).

Example 45

(±)-4-(8-Azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2543)

[0417]



[0418] In a similar manner to the procedures described in Examples 9(i), 9(ii) and 9(iii) above, coupling, debenzyla-
tion, dehydration and desilylation reactions were carried out using 8-benzyl-8-azabicyclo[3.2.1]octane-3-one as a start-
ing material instead of 1-benzylpiperidine-4-one to give the title compound (yield 35%) as a pale yellow powder.

[0419] Melting point: 212 - 213°C

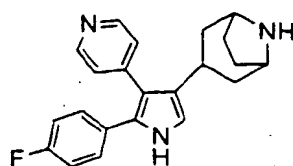
[0420] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.32 (1H, broad singlet);
8.45 (2H, doublet, J=6 Hz);
7.17-7.07 (6H, multiplet);
6.31 (1H, doublet, J=3 Hz);
5.45 (1H, doublet, J=6 Hz);
3.53-3.46 (1H, multiplet);
3.38-3.32 (1H, multiplet);
2.55-2.48 (1H, broad multiplet);
1.36-1.69 (3H, multiplet);
1.66-1.55 (1H, multiplet);
1.52-1.44 (1H, multiplet).

Example 46

4-(8-Azabicyclo[3.2.1]octan-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2541)

[0421]



[0422] In a similar manner to that described in Example 11 above, (±)-4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 45 above] was reduced to give the title compound (yield 74%) as a white powder.

[0423] Melting point: 235 - 237°C (decomposition)

[0424] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

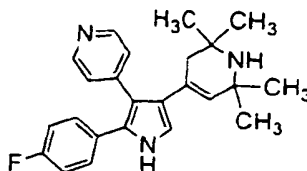
11.14 (1H, broad singlet);
8.48 (2H, doublet, J=6 Hz);
7.15-7.05 (6H, multiplet);
6.73 (1H, doublet, J=2 Hz);
3.18-3.15 (2H, multiplet);
2.36-2.78 (1H, multiplet);
2.05-1.96 (2H, multiplet);
1.63-1.56 (2H, multiplet);

1.50-1.43 (2H, multiplet);
1.17-1.08 (2H, multiplet).

Example 47

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydro-2,2,6,6-tetramethylpyridin-4-yl)-1H-pyrrole (Compound No. 1-2650)

[0425]



[0426] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 2,2,6,6-tetramethylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 50%) as a white powder.

[0427] Melting point: 258 - 259°C

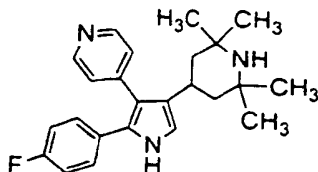
[0428] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.37 (1H, broad singlet);
8.42 (2H, doublet, J=6 Hz);
7.19-7.09 (6H, multiplet);
6.36 (1H, doublet, J=2 Hz);
5.04-5.02 (1H, broad multiplet);
1.93 (2H, singlet);
1.25-1.15 (1H, broad singlet);
1.04 (6H, singlet);
0.94 (6H, singlet).

Example 48

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(2,2,6,6-tetramethylpiperidin-4-yl)-1H-pyrrole (Compound No. 1-2523)

[0429]



[0430] In a similar manner to that described in Example 11 above, 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydro-2,2,6,6-tetramethylpyridin-4-yl)-1H-pyrrole [prepared as described in Example 47 above] was reduced to give the title compound (yield 77%) as a white powder.

[0431] Melting point: 242 - 243°C

[0432] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

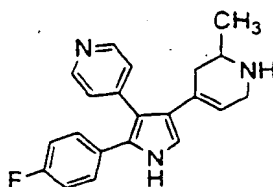
11.18 (1H, broad singlet);
8.47 (2H, doublet, J=6 Hz);
7.21-7.15 (4H, multiplet);
7.11 (2H, triplet, J=9 Hz);

6.70 (1H, doublet, J=2 Hz);
 2.95 (1H, triplet of triplets, J=13 Hz, 3 Hz);
 1.50 (2H, doublet of doublets, J=13 Hz, 3 Hz);
 1.09 (2H, triplet, J=13 Hz);
 1.01 (6H, singlet);
 0.96 (6H, singlet).

Example 49

(±)-2-(4-Fluorophenyl)-4-(2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2707)

[0433]



[0434] In a similar manner to the procedures described in Examples 9(i), 9(ii) and 9(iii) above, coupling, debenzyla-
 tion, dehydration and desilylation reactions were carried out using (±)-1-benzyl-2-methylpiperidine-4-one as a starting
 material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (6-methyl-1,2,3,6-tet-
 rahydropyridin-4-yl) isomer thereof (which is the title compound of Example 50). The mixture was separated by chro-
 matography on a silica gel column using a 25:1 by volume mixture of ethyl acetate and isopropylamine as the eluant
 to give the title compound (yield 3%, R_f value= 0.50) as a white powder.

[0435] Melting point: 178 - 180°C (decomposition)

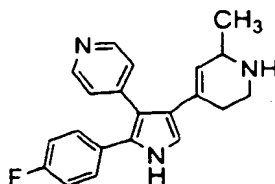
[0436] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

8.36 (2H, doublet, J=6 Hz);
 7.24 (2H, doublet, J=6 Hz);
 7.17 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 7.00 (2H, triplet, J=9 Hz);
 6.85 (1H, singlet);
 5.40-5.35 (1H, broad multiplet);
 3.46-3.38 (1H, multiplet);
 3.36-3.28 (1H, multiplet);
 2.96-2.88 (1H, multiplet);
 2.23-2.15 (1H, multiplet);
 2.06-1.97 (1H, multiplet);
 1.14 (3H, doublet, J=6 Hz).

Example 50

(±)-2-(4-Fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2665)

[0437]



[0438] The title compound (yield 5%, Rf value = 0.45) was obtained as a white powder during the chromatography performed in Example 49 above.

[0439] Melting point: 201 - 204°C (decomposition)

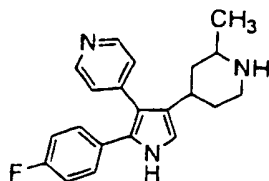
[0440] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

- 5 8.36 (2H, doublet, J=6 Hz);
 7.24 (2H, doublet, J=6 Hz);
 7.18 (2H, doublet of doublets, J=9 Hz, 6 Hz);
 7.00 (2H, triplet, J=9 Hz);
 6.36 (1H, singlet);
 10 5.24-5.20 (1H, multiplet);
 3.48-3.40 (1H, multiplet);
 3.16-3.09 (1H, multiplet);
 2.90-2.82 (1H, multiplet);
 2.37-2.27 (1H, multiplet);
 15 2.18-2.10 (1H, multiplet);
 1.05 (3H, doublet, J=7 Hz).

Example 51

20 (±)-2-(4-Fluorophenyl)-4-(2-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2522)

[0441]



35 [0442] In a similar manner to the procedure in Example 11 above, (±)-2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 50 above] was reduced to give the title compound (yield 83%) as a pale yellow powder.

[0443] Melting point: 198 - 205°C

[0444] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

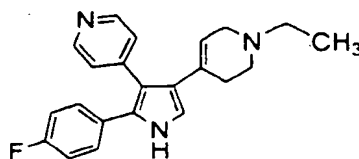
- 40 8.53 (2H, doublet, J=4 Hz);
 8.33 (0.5H, broad singlet);
 8.31 (0.5H, broad singlet);
 7.14 (2H, doublet, J=4 Hz);
 7.11 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.95 (2H, triplet, J=9 Hz);
 45 6.79 (0.5H, doublet, J=3 Hz);
 6.74 (0.5H, doublet, J=3 Hz);
 3.25-2.63 (4H, multiplet);
 1.74-1.43 (5H, multiplet);
 50 1.11 (1.5H, doublet, J=7 Hz);
 1.07 (1.5H, doublet, J=7 Hz).

55

Example 52

4-(1-Ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2629)

[0445]



[0446] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-ethylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 55%) as a pale yellow powder.

[0447] Melting point: 234 - 236°C (decomposition)

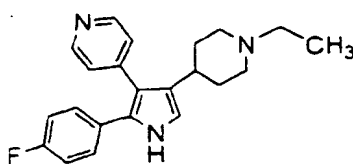
[0448] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.36 (1H, broad singlet);
8.45 (2H, doublet, J=6 Hz);
7.18-7.08 (6H, multiplet);
6.90 (1H, doublet, J=3 Hz);
5.26-5.22 (1H, broad multiplet);
2.34-2.79 (2H, broad multiplet);
2.45 (2H, triplet, J=6 Hz);
2.34 (2H, quartet, J=8 Hz);
2.18-2.12 (2H, broad multiplet);
0.99 (3H, triplet, J=8 Hz).

Example 53

4-(1-Ethylpiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-146)

[0449]



[0450] In a similar manner to the procedure described in Example 11 above, 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in Example 52 above] was reduced to give the title compound (yield 95%) as a white powder.

[0451] Melting point: 262 - 263°C (decomposition)

[0452] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

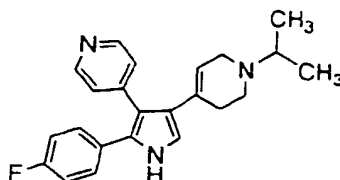
11.19 (1H, broad singlet);
8.48 (2H, doublet, J=6 Hz);
7.18-7.07 (6H, multiplet);
6.76 (1H, doublet, J=2 Hz);
2.87-2.79 (2H, broad multiplet);
2.37 (1H, triplet of triplets, J=12 Hz, 4 Hz);
2.26 (2H, quartet, J=7 Hz);
1.80-1.71 (2H, broad multiplet);

1.66-1.59 (2H, broad multiplet);
 1.55-1.43 (2H, multiplet);
 0.96 (3H, triplet, J=7 Hz).

5 Example 54

2-(4-Fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2632)

10 [0453]



20 [0454] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-isopropylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 42%) as a pale yellow powder.

[0455] Melting point: 237 - 240°C (decomposition)

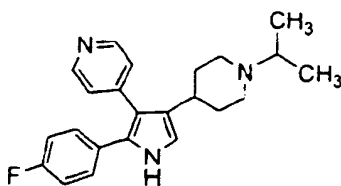
[0456] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

25 11.35 (1H, broad singlet);
 8.45 (2H, doublet, J=6 Hz);
 7.18-7.08 (6H, multiplet);
 6.38 (1H, doublet, J=3 Hz);
 5.28-5.24 (1H, broad multiplet);
 30 2.94-2.89 (2H, broad multiplet);
 2.64 (1H, septet, J=6 Hz);
 2.49 (2H, triplet, J=6 Hz);
 2.15-2.09 (2H, broad multiplet);
 0.96 (6H, doublet, J=6 Hz).

35 Example 55

2-(4-Fluorophenyl)-4-(1-isopropylpiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2633)

40 [0457]



50 [0458] In a similar manner to that described in Example 11 above, 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 54 above) was reduced to give the title compound (yield 80%) as white needle-like crystals.

[0459] Melting point: 264 - 266°C (decomposition)

55 [0460] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

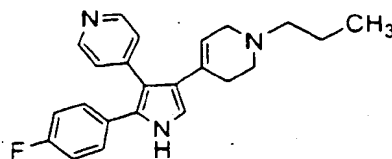
11.18 (1H, broad singlet);
 8.48 (2H, doublet, J=6 Hz);
 7.17-7.06 (6H, multiplet);

6.76 (1H, singlet);
 2.77-2.70 (2H, broad multiplet);
 2.63 (1H, septet, J=6 Hz);
 2.35 (1H, triplet of triplets, J=12 Hz, 4 Hz);
 2.06-1.97 (2H, multiplet);
 1.67-1.60 (2H, broad multiplet);
 1.52-1.40 (2H, multiplet);
 0.92 (6H, doublet, J=6 Hz).

Example 56

2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2630)

[0461]



[0462] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-propylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 46%) as a white powder.

[0463] Melting point: 236 - 238°C (decomposition)

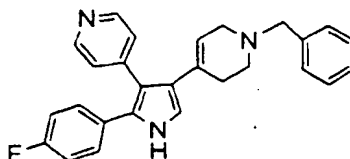
[0464] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.36 (1H, broad singlet);
 8.45 (2H, doublet, J=6 Hz);
 7.18-7.08 (5H, multiplet);
 6.90 (1H, doublet, J=3 Hz);
 5.26-5.23 (1H, multiplet);
 2.84-2.79 (2H, multiplet);
 2.44 (2H, triplet, J=5 Hz);
 2.25 (2H, triplet, J=8 Hz);
 2.17-2.11 (2H, broad multiplet);
 1.43 (2H, triplet of quartets, J=8 Hz, 8 Hz);
 0.84 (3H, triplet, J=8 Hz).

Example 57

4-(1-Benzyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2644)

[0465]



[0466] In a similar manner to that described in Example 9(iii) above, dehydration and desilylation reactions were carried out using 4-(1-benzyl-4-hydroxypiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in Example 9(i) above] to give the title compound (yield 88%) as a pale yellow powder.

[0467] Melting point: 217 - 219°C (decomposition)

[0468] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.37 (1H, broad singlet);

8.45 (2H, doublet, J=6 Hz);

7.34-7.21 (5H, multiplet);

7.18-7.08 (5H, multiplet);

6.90 (1H, doublet, J=3 Hz);

5.23-5.21 (1H, broad multiplet);

3.50 (2H, singlet);

2.36-2.81 (2H, broad multiplet)

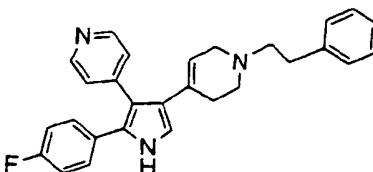
2.48 (2H, triplet, J=6 Hz);

2.18-2.12 (2H, broad multiplet).

Example 58

2-(4-Fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2645)

[0469]



[0470] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-phenethylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 46%) as a pale brown powder.

[0471] Melting point: 219 - 221°C (decomposition)

[0472] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.37 (1H, broad singlet);

8.45 (2H, doublet, J=6 Hz);

7.29-7.09 (11H, multiplet);

6.90 (1H, doublet, J=3 Hz);

5.28-5.24 (1H, broad multiplet);

2.95-2.89 (2H, broad multiplet);

2.76-2.69 (2H, multiplet);

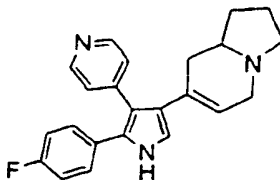
2.57-2.51 (4H, multiplet);

2.19-2.12 (2H, broad multiplet).

Example 59

(±)-2-(4-Fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2654)

[0473]



[0474] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and

desilylation reactions were carried out using (\pm)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one as a starting material instead of 1-benzylpiperidine-4-one to give a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizine-7-yl) isomer thereof (which is the compound of Example 60 below). The mixture was separated by chromatography on a silica gel column using a 10:1:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give the title compound (yield 24%, Rf value = 0.50) as a white powder.

[0475] Melting point: 220 - 222°C (decomposition)

[0476] ^1H -Nuclear magnetic resonance spectrum (500 MHz, CD_3OD) δ ppm:

8.36 (2H, doublet, J=6 Hz);

7.24 (2H, doublet, J=6 Hz);

7.17 (2H, doublet of doublets, J=9 Hz, 6 Hz);

7.00 (2H, triplet, J=9 Hz);

6.35 (1H, singlet);

5.41-5.37 (1H, multiplet);

3.47-3.40 (1H, multiplet);

3.18-3.12 (1H, multiplet);

2.34-2.77 (1H, multiplet);

2.40-2.26 (2H, multiplet);

2.23-2.11 (2H, multiplet);

2.05-1.97 (1H, multiplet);

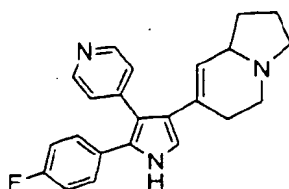
1.90-1.75 (2H, multiplet);

1.46-1.36 (1H, multiplet).

Example 60

(\pm)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2653)

[0477]



[0478] The title compound (yield 27%, Rf value = 0.40) was obtained as a pale yellow powder during the chromatography performed in Example 59 above.

[0479] Melting point: 188 - 190°C (decomposition)

[0480] ^1H -Nuclear magnetic resonance spectrum (500 MHz, CD_3OD) δ ppm:

8.37 (2H, doublet, J=6 Hz);

7.23 (2H, doublet, J=6 Hz);

7.18 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.00 (2H, triplet, J=9 Hz);

6.36 (1H, singlet);

5.34-5.30 (1H, broad multiplet);

3.38-3.30 (1H, multiplet);

2.99-2.86 (2H, multiplet);

2.32-2.73 (2H, multiplet);

2.40-2.31 (1H, multiplet);

2.25-2.16 (1H, multiplet);

1.97-1.86 (2H, multiplet);

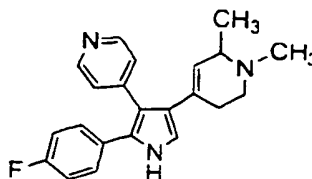
1.84-1.73 (1H, multiplet);

1.48-1.38 (1H, multiplet).

Example 61

(±)-4-(1,6-Dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2666)

[0481]



[0482] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-1,2-dimethylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give a mixture of the title compound and the (1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl) isomer thereof (which is the compound of Example 62 below). The mixture was separated by chromatography on a silica gel column using a 99:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 10%, Rf value = 0.40) as a pale yellow powder.

[0483] Melting point: 205 - 209°C (decomposition)

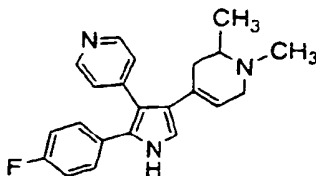
[0484] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
 8.29-8.10 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);
 7.13 (2H, doublet of doublets J=9 Hz, 5 Hz);
 6.97 (2H, triplet, J=9 Hz);
 6.33 (1H, doublet, J=3 Hz);
 5.19 (1H, singlet);
 2.90-2.83 (1H, multiplet);
 2.55-2.32 (2H, multiplet);
 2.37 (3H, singlet);
 2.17-2.07 (2H, multiplet);
 1.05 (3H, doublet, J=7 Hz).

Example 62

(±)-4-(1,2-Dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2708)

[0485]



[0486] The title compound (yield 14%, Rf value = 0.30) was obtained as a pale yellow powder during the chromatography performed in Example 61 above.

[0487] Melting point: 198 - 201 °C

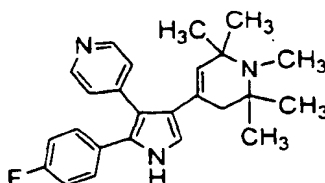
[0488] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:
 8.47 (2H, doublet, J=6 Hz);

8.33 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);
 7.12 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.97 (2H, triplet, J=9 Hz);
 6.82 (1H, doublet, J=3 Hz);
 5.40 (1H, singlet);
 3.22-3.18 (1H, multiplet);
 2.95-2.91 (1H, multiplet);
 2.57-2.46 (1H, multiplet);
 2.34 (3H, singlet);
 2.27-2.23 (1H, multiplet);
 2.11-2.09 (1H, multiplet);
 1.08 (3H, doublet, J=6 Hz).

Example 63

2-(4-Fluorophenyl)-4-(1,2,2,6,6-pentamethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2651)

[0489]



[0490] In a similar manner to the procedures described in Example 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1,2,2,6,6-pentamethylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 69%) as a white powder.

[0491] Melting point: 264 - 268°C (decomposition)

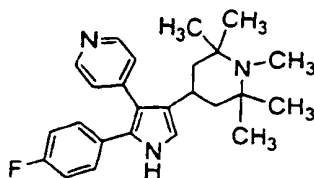
[0492] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.38 (1H, broad singlet);
 8.43 (2H, doublet, J=6 Hz);
 7.20-7.09 (6H, multiplet);
 6.88 (1H, doublet, J=3 Hz);
 4.88 (1H, singlet);
 2.18 (3H, singlet);
 2.07 (2H, singlet);
 1.01 (6H, singlet);
 0.90 (6H, singlet).

Example 64

2-(4-Fluorophenyl)-4-(1,2,2,6,6-pentamethylpiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2652)

[0493]



[0494] In a similar manner to that described in Example 11 above, 2-(4-fluorophenyl)-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 63 above) was reduced to give the title compound (yield 82%) as a white powder.

[0495] Melting point: 248 - 250°C (decomposition)

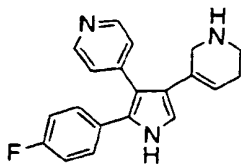
[0496] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.19 (1H, broad singlet);
8.47 (2H, doublet, J=6 Hz);
7.21-7.15 (4H, multiplet);
7.11 (2H, triplet, J=9 Hz);
6.73 (1H, doublet, J=2 Hz);
2.91-2.62 (1H, multiplet);
2.14 (3H, singlet);
1.54-1.47 (2H, multiplet);
1.36 (2H, triplet, J=13 Hz);
1.01 (6H, singlet);
0.36 (6H, singlet).

Example 65

2-(4-Fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-5-yl)-1H-pyrrole (Compound No. 1-2627)

[0497]



[0498] In a similar manner to the procedures described in Example 9(i), 9(ii) and 9(iii) above, coupling, debenzylation, dehydration and desilylation reactions were carried out using 1-benzylpiperidine-3-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 12%) as a pale brown powder.

[0499] Melting point: 230 - 233°C (decomposition)

[0500] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CD₃OD) δ ppm:

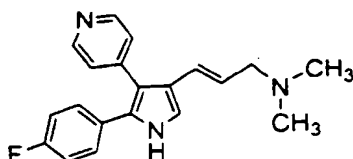
8.36 (2H, doublet, J=6 Hz);
7.24 (2H, doublet, J=6 Hz);
7.18 (2H, doublet of doublets, J=9 Hz, 6 Hz);
7.01 (2H, triplet, J=9 Hz);
6.32 (1H, singlet);
5.55-5.50 (1H, broad multiplet);

3.36-3.30 (2H, multiplet);
 2.90 (2H, triplet, J=6 Hz);
 2.16-2.09 (2H, broad multiplet).

Example 66

4-(3-Dimethylamino-1-propen-1-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2519)

[0501]



66(i) 2-(4-Fluorophenyl)-4-(3-hydroxy-1-propen-1-yl)-3-(pyridin-4-yl)-1H-pyrrole

[0502] 27.6 ml (27.6 mmol) of a 1M solution of diisobutylaluminum hydride in dichloromethane were added dropwise to a solution of 3.88 g (12.5 mmol) of ethyl 3-[2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrol-4-yl]acrylate in 120 ml of dichloromethane at -78°C with stirring. The resulting mixture was stirred at the same temperature for 3 hours. At the end of this time, the reaction mixture was added to a saturated aqueous solution of sodium hydrogencarbonate and the resulting mixture was filtered. The filtrate was extracted with dichloromethane and the organic extract was washed with water, dried over anhydrous sodium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 9:1 by volume mixture of hexane and ethyl acetate as the eluant to afford 3.42 g (yield 93%) of the title compound as a pale brown powder.

[0503] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CD₃OD) δ ppm:

8.41 (2H, doublet, J=6 Hz);
 7.23 (2H, doublet, J=6 Hz);
 7.20 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.08 (1H, singlet);
 7.01 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet, J=16 Hz);
 5.98 (1H, doublet of triplets, J=16 Hz, 6 Hz);
 4.09 (2H, doublet of doublets, J=6 Hz, 1 Hz).

66(ii) 3-[2-(4-Fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrol-4-yl]acrylaldehyde

[0504] 16.32 g of activated manganese oxide were added to a solution of 1.63 g (5.54 mmol) of 2-(4-fluorophenyl)-4-(3-hydroxy-1-propen-1-yl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 66(i) above] in 80 ml of dimethylsulfoxide. The resulting mixture was stirred at room temperature for 24 hours. At the end of this time, ethyl acetate was added to the reaction mixture and then this was filtered. Water was added to the resulting filtrate and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was washed with dichloromethane to afford 1.36 g (yield 84%) of the title compound as a pale brown powder.

[0505] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CD₃OD-CDCl₃) δ ppm:

9.42 (1H, doublet, J=8 Hz);
 8.51 (2H, doublet, J=6 Hz);
 7.49 (1H, singlet);
 7.45 (1H, doublet, J=16 Hz);
 7.27 (2H, doublet, J=6 Hz);
 7.22 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.01 (2H, triplet, J=9 Hz);
 6.39 (1H, doublet of doublets, J=16 Hz, 8 Hz).

66(iii) 4-(3-Dimethylamino-1-propen-1-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0506] 700 mg (2.38 mmol) of 3-[2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrol-4-yl]acrylaldehyde [prepared as described in step 66 (i) above] and 1.95 g (23.8 mmol) of dimethylamine hydrochloride were dissolved in 110 ml of a 1:1 by volume mixture of methanol and tetrahydrofuran, after which 226 mg (3.59 mmol) of sodium cyanotrihydridoborate were added to the solution. The resulting mixture was stirred at room temperature for 19 hours. At the end of this time, the reaction mixture was added to a saturated aqueous solution of sodium hydrogencarbonate and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 20:1:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give 429 mg (yield 56%) of the title compound as a white powder.

[0507] Melting point: 196 - 198°C (decomposition)

[0508] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CD₃OD-CDCl₃) δ ppm:

8.42 (2H, doublet, J=6 Hz);

7.23 (2H, doublet, J=6 Hz);

7.20 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.09 (1H, singlet);

6.98 (2H, triplet, J=9 Hz);

6.35 (1H, doublet, J=16 Hz);

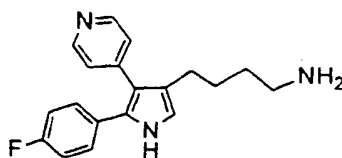
5.38 (1H, doublet of triplets, J=16 Hz, 7 Hz);

3.04 (2H, doublet, J=7 Hz);

2.27 (6H, singlet).

Example 674-(4-Aminobutyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-98)

[0509]

67(i) 4-(3-Cyanopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0510] 568 mg (8.72 mmol) of potassium cyanide were added to a solution of 3.04 g (7.26 mmol) of 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-[3-(p-toluene-sulfonyloxy)propyl]-1H-pyrrole [prepared as described in Example 23(iv) above] in 60 ml of N,N-dimethylformamide and the resulting mixture was stirred at 100°C for 2 hours. After cooling to room temperature, a saturated aqueous solution of sodium hydrogencarbonate was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 2:1 by volume mixture of hexane and ethyl acetate as the eluant to give 0.94 g (yield 42%) of the title compound as a pale yellow powder.

[0511] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.53 (2H, doublet, J=6 Hz);

8.25 (1H, broad singlet);

7.13 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.12 (2H, doublet, J=6 Hz);

6.97 (2H, triplet, J=9 Hz);

6.75 (1H, doublet, J=3 Hz);

2.70 (2H, triplet, J=7 Hz);

2.28 (2H, triplet, J=7 Hz);

1.76 (2H, quintet, J=7 Hz).

67(ii) 4-(4-Aminobutyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole

[0512] A solution of 0.94 g (3.08 mmol) of 4-(3-cyanopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole [prepared as described in step 67(i) above] in 50 ml of tetrahydrofuran was added dropwise to a suspension of 468 mg (12.32 mmol) of lithium aluminum hydride in 100 ml of tetrahydrofuran at room temperature with stirring. The resulting mixture was stirred at 60°C for 3 hours. After cooling to room temperature, 0.47 ml of a 15% aqueous solution of sodium hydroxide and 20 ml of water were successively added gradually to the reaction mixture. The resulting mixture was filtered and the filtrate thus obtained was concentrated by evaporation under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using a 25:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give 410 mg (yield 43%) of the title compound as a pale red powder.

[0513] Melting point: 185 - 187°C

[0514] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

10.02 (1H, broad singlet);

8.49 (2H, doublet, J=4 Hz);

7.19 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.14 (2H, doublet, J=4 Hz);

6.94 (2H, triplet, J=9 Hz);

6.72 (1H, doublet, J=2 Hz);

2.65 (2H, triplet, J=7 Hz);

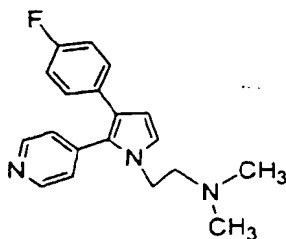
2.51 (2H, doublet, J=7 Hz);

2.46-2.08 (2H, multiplet);

1.52-1.46 (4H, multiplet).

Example 681-(2-Dimethylaminoethyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-106)

[0515]



[0516] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 2-dimethylaminoethylamine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (yield 8%) as a pale brown powder.

[0517] Melting point: 70-73°C

[0518] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.60 (2H, doublet, J=6 Hz);

7.19 (2H, doublet, J=6 Hz);

7.07 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.39 (2H, triplet, J=9 Hz);

6.38 (1H, doublet, J=3 Hz);

6.36 (1H, doublet, J=3 Hz);

3.96 (2H, triplet, J=7 Hz);

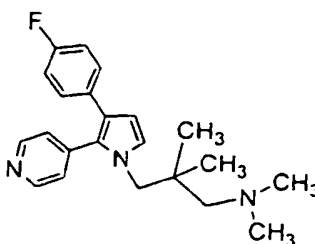
2.48 (2H, triplet, J=7 Hz);

2.15 (6H, singlet).

Example 69

1-(2,2-Dimethyl-3-dimethylaminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2626)

[0519]



[0520] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 2,2-dimethyl-3-dimethylaminopropylamine as a starting material instead of 4-amino-1-benzylpiperidine to give the title compound (yield 21%) as a white powder.

[0521] Melting point: 101 - 103°C

[0522] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.58 (2H, doublet, J=6 Hz);

7.14 (2H, doublet, J=6 Hz);

7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.38 (2H, triplet, J=9 Hz);

6.34 (1H, doublet, J=3 Hz);

6.32 (1H, doublet, J=3 Hz);

3.90 (2H, singlet);

2.22 (6H, singlet);

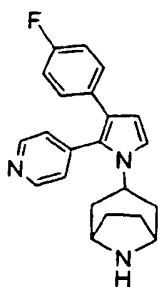
2.00 (2H, singlet);

0.64 (6H, singlet).

Example 70

1-(8-Azabicyclo[3.2.1]octan-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2541)

[0523]



[0524] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 3-amino-3-benzyl-8-azabicyclo[3.2.1]octane as a starting material instead of 4-amino-1-benzylpiperidine to afford the N-benzyl derivative of the title compound. The N-benzyl derivative was then debenzylated in a similar manner to that described in Example 9(ii) above to give the title compound (yield 90%) as a white powder.

[0525] Melting point: 204- 205°C

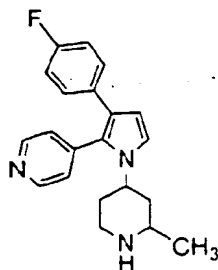
[0526] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.60 (2H, doublet, J=6 Hz);
 7.14 (2H, doublet, J=6 Hz);
 7.02 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.38 (1H, doublet, J=3 Hz);
 5 6.36 (2H, triplet, J=9 Hz);
 6.38 (1H, doublet, J=3 Hz);
 4.36-4.25 (1H, multiplet);
 3.67-3.53 (2H, multiplet);
 2.51-2.39 (2H, multiplet);
 10 1.90-1.67 (4H, multiplet);
 1.57-1.47 (2H, multiplet).

Example 71

15 (±)-3-(4-Fluorophenyl)-1-(2-methylpiperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole (Compound No. 2-2522)

[0527]



20 [0528] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using (±)-4-amino-1-benzyl-2-methylpiperidine as a starting material instead of 4-amino-1-benzylpiperidine to afford the N-benzyl derivative of the title compound. The N-benzyl derivative was then debenzylated in a similar manner to that described in Example 9(ii) above to give the title compound (yield 24%) as a white powder.

35 [0529] Melting point: 173 - 175°C

[0530] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

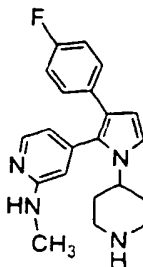
8.61 (2H, doublet, J=5 Hz);
 7.15 (1H, doublet, J=5 Hz);
 7.14 (1H, doublet, J=5 Hz);
 40 7.07-7.03 (2H, multiplet);
 7.00 (0.5H, doublet, J=3 Hz);
 6.95 (0.5H, doublet, J=3 Hz);
 6.38 (2H, triplet, J=9 Hz);
 6.38 (1H, doublet, J=3 Hz);
 45 4.30-4.26 (0.5H, multiplet);
 4.00-3.91 (0.5H, multiplet);
 3.47-3.39 (0.5H, multiplet);
 3.17-3.15 (0.5H, multiplet);
 2.98-2.88 (1H, multiplet);
 50 2.66-2.61 (1H, multiplet);
 2.06-1.76 (4H, multiplet);
 1.11 (1.5H, doublet, J=7 Hz);
 1.08 (1.5H, doublet, J=7 Hz).

Example 723-(4-Fluorophenyl)-2-(2-methylaminopyridin-4-yl)-1-(piperidin-4-yl)-1H-pyrrole (Compound No. 2-2042)

5 [0531]

10

15

20 72(i) 1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-fluoropyridin-4-yl)-1H-pyrrole

[0532] In a similar manner to the procedures described in Examples 12(i), 12(ii) and 12(iii) above, reactions were carried out using 2-fluoropyridin-4-carboxyaldehyde as a starting material instead of pyridine-4-carboxyaldehyde to give the title compound (yield 49%) as a pale brown powder.

25 [0533] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.20 (1H, doublet, J=5 Hz);

7.37-7.24 (5H, multiplet);

7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);

7.02-6.96 (2H, multiplet);

30 6.90 (2H, triplet, J=9 Hz);

6.75 (1H, singlet);

6.37 (1H, doublet, J=3 Hz);

3.89-3.78 (1H, multiplet);

3.52 (2H, singlet);

35 3.01-2.96 (2H, multiplet);

2.12-1.96 (4H, multiplet);

1.94-1.86 (2H, multiplet).

40 72(ii) 1-(1-Benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylaminopyridin-4-yl)-1H-pyrrole

[0534] A mixture of a solution of 1.50 g (3.49 mmol) of 1-(1-benzylpiperidin-4-yl)-3-(4-fluorophenyl)-2-(2-fluoropyridin-4-yl)-1H-pyrrole [prepared as described in step 72(i) above] in 30 ml of tetrahydrofuran and 50 ml (100 mmol) of a 2M solution of methylamine in tetrahydrofuran was stirred at 150°C in an autoclave for 28 hours. At the end of this time, the solvent was distilled off from the reaction mixture under reduced pressure. Ethyl acetate and water were added to the residue, and the organic layer was separated, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 20:1 by volume mixture of dichloromethane and methanol as the eluant to give 1.83 g (yield 60%) of the title compound as a pale brown amorphous solid.

50 [0535] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.10 (1H, doublet, J=5 Hz);

7.35-7.23 (5H, multiplet);

7.12 (2H, doublet of doublets, J=9 Hz, 5 Hz);

6.92 (1H, doublet, J=3 Hz);

6.88 (2H, triplet, J=9 Hz);

55 6.48 (1H, doublet, J=5 Hz);

6.37 (1H, doublet, J=3 Hz);

6.20 (1H, singlet);

4.66-4.57 (1H, multiplet);

3.93-3.84 (1H, multiplet);
 3.51 (2H, singlet);
 2.99-2.91 (2H, multiplet);
 2.34 (3H, doublet, J=5 Hz);
 2.06-1.85 (6H, multiplet).

72(iii) 3-(4-Fluorophenyl)-2-(2-methylaminopyridin-4-yl)-1-(piperidin-4-yl)-1H-pyrrole

[0536] In a similar manner to that described in Example 9(ii) above, 1-(1-benzyl-piperidin-4-yl)-3-(4-fluorophenyl)-2-(2-methylaminopyridin-4-yl)-1H-pyrrole [prepared as described in step 72(ii) above] was debenzylated to give the title compound (yield 52%) as a white powder.

[0537] Melting point: 165 - 167°C

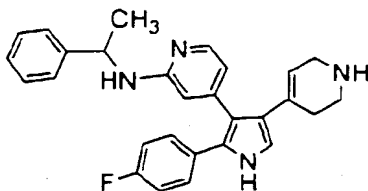
[0538] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.12 (1H, doublet, J=5 Hz);
 7.13 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.93 (1H, doublet, J=3 Hz);
 6.38 (2H, triplet, J=9 Hz);
 6.49 (1H, doublet of doublets, J=5 Hz, 1 Hz);
 6.38 (1H, doublet, J=3 Hz);
 6.21 (1H, doublet, J=1 Hz);
 4.65-4.55 (1H, multiplet);
 4.02-3.97 (1H, multiplet);
 3.16 (2H, broad doublet, J=12 Hz);
 2.35 (3H, doublet, J=5 Hz);
 2.60 (2H, doublet of triplets, J=12 Hz, 3 Hz);
 1.99-1.81 (4H, multiplet).

Example 73

2-(4-Fluorophenyl)-3-[2-(1(S)-phenylethylamino)pyridin-4-yl]-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole
 (Compound No. 1-2600)

[0539]



73(i) 4-Bromo-2-(4-fluorophenyl)-3-(2-fluoropyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0540] In a similar manner to the procedures described in Examples 1(i), 7(i), 7(ii) and 7(iii) above, a pyrrole ring-forming reaction, a decarboxylation reaction, a silylation reaction and a bromination reaction were carried out using ethyl 3-(2-fluoropyridin-4-yl)acrylate as a starting material instead of ethyl 3-(4-pyridyl)acrylate to give the title compound (yield 48%) as a pale yellow powder.

[0541] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

8.05 (1H, doublet, J=5 Hz);
 7.38 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.20 (2H, triplet, J=9 Hz);
 7.15 (1H, singlet);
 7.01-6.97 (1H, multiplet);
 6.78 (1H, singlet);
 1.20-1.07 (3H, multiplet);
 0.99 (18H, doublet, J=7 Hz).

73(ii) 4-(1-Allyloxycarbonyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(2-fluoropyridin-4-yl)-1H-pyrrole

[0542] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-(4-fluorophenyl)-3-(2-fluoropyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 73(i) above] instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole and using 1-allyloxycarbonylpiperidine-4-one instead of 1-benzylpiperidine-4-one to give the title compound (yield 6%) as a white powder.

[0543] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.42-8.31 (1H, broad singlet);
 8.09 (1H, doublet, J=5 Hz);
 7.15 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 7.04-6.97 (3H, multiplet);
 6.32 (1H, singlet);
 6.77 (1H, singlet);
 6.00-5.89 (1H, multiplet);
 5.63-5.39 (1H, multiplet);
 5.31 (1H, doublet, J=16 Hz);
 5.21 (1H, doublet, J=11 Hz);
 4.62 (2H, doublet, J=5 Hz);
 4.04-3.96 (2H, broad singlet);
 3.63-3.52 (2H, broad singlet);
 2.29-2.12 (2H, multiplet).

73(iii) 4-(1-Allyloxycarbonyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-[2-(1(S)-phenylethylamino)pyridin-4-yl]-1H-pyrrole

[0544] A mixture of 360 mg (0.85 mmol) of 4-(1-allyloxycarbonyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(2-fluoropyridin-4-yl)-1H-pyrrole [prepared as described in step 73 (ii) above], 3.6 ml (28.3 mmol) of 1(S)-phenylethylamine and 0.36 ml of concentrated hydrochloric acid was stirred at 150°C for 10 hours. At the end of this time, a saturated aqueous solution of sodium hydrogencarbonate was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water, dried over anhydrous magnesium sulfate and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 1:1 by volume mixture of hexane and ethyl acetate as the eluant to give 229 mg (yield 49%) of the title compound as a brown amorphous solid.

[0545] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.20-8.12 (1H, broad singlet);
 7.96 (1H, doublet, J=5 Hz);
 7.25 (4H, singlet);
 7.22-7.16 (1H, multiplet);
 7.04 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.39 (2H, triplet, J=9 Hz);
 6.74 (1H, singlet);
 6.44 (1H, doublet, J=4 Hz);
 6.10 (1H, singlet);
 6.00-5.89 (1H, multiplet);
 5.54-5.33 (1H, multiplet);
 5.32 (1H, doublet, J=16 Hz);
 5.22 (1H, doublet, J=11 Hz);
 5.05 (1H, broad doublet, J=6 Hz);
 4.62 (2H, doublet, J=5 Hz);
 4.54-4.46 (1H, multiplet);
 3.93-3.83 (2H, multiplet);
 3.50-3.38 (2H, multiplet);
 2.20-2.05 (2H, multiplet);
 1.48 (3H, doublet, J=7 Hz).

73(iv) 2-(4-Fluorophenyl)-3-[2-(1(S)-phenylethylamino)pyridin-4-yl]-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole

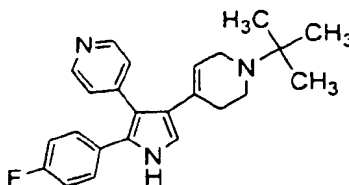
[0546] A mixture of 229 mg (0.44 mmol) of 4-(1-allyloxycarbonyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-[2-(1(S)-phenylethylamino)pyridin-4-yl]-1H-pyrrole [prepared as described in step 73(iii) above] in 2.5 ml of a 10:1 by volume mixture of dioxane and water, 55 μ l (0.66 mmol) of pyrrolidine and 5 mg (0.0044 mmol) of tetrakis(triphenylphosphine)palladium (0) was stirred at 0°C for 10 minutes. At the end of this time, water was added to the reaction mixture and then this was extracted with ethyl acetate. The organic extract was washed with water and then concentrated by evaporation under reduced pressure. The residue thus obtained was purified by chromatography on a silica gel column using a 10:1:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give 41 mg (yield 21%) of the title compound as a pale brown amorphous solid.

[0547] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.19-8.01 (1H broad singlet);
 7.96 (1H doublet J=5 Hz);
 7.26 (4H singlet);
 7.24-7.15 (1H multiplet);
 7.06 (2H doublet of doublets, J=9 Hz, 5 Hz);
 6.89 (2H triplet J=9 Hz);
 6.74 (1H singlet);
 6.47 (1H doublet J=4 Hz);
 6.15 (1H singlet);
 5.47 (1H doublet J=3 Hz);
 5.05 (1H broad doublet, J=6 Hz);
 4.60-4.52 (1H multiplet);
 3.31-3.26 (2H multiplet);
 2.94-2.84 (2H multiplet);
 2.13-2.07 (2H multiplet);
 1.48 (3H doublet J=7 Hz).

Example 744-(1-t-Butyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2635)

[0548]



[0549] In a similar manner to the procedures described in Examples 9(i), 9(ii) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-t-butylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 20%) as a pale yellow powder.

[0550] Melting point: 242 - 244°C

[0551] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃-DMSO-d₆) δ ppm:

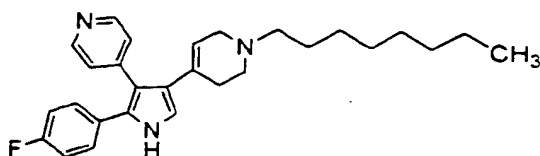
9.72 (1H, broad singlet);
 8.46 (2H, doublet, J=6 Hz);
 7.19 (2H, doublet, J=6 Hz);
 7.16 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.94 (2H, triplet, J=9 Hz);
 6.81 (1H, doublet, J=3 Hz);
 5.46 (1H, singlet);
 3.19-3.11 (2H, multiplet);
 2.66-2.58 (2H, multiplet);

2.26-2.18 (2H, multiplet);
1.10 (9H, singlet).

Example 75

2-(4-Fluorophenyl)-4-(1-octyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2639)

[0552]



[0553] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 1-octylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to give the title compound (yield 26%) as a white powder.

[0554] Melting point: 173 - 175°C

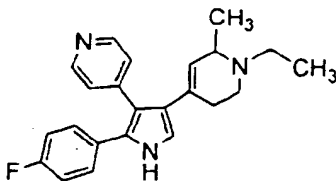
[0555] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
8.29 (1H, broad singlet);
7.17 (2H, doublet, J=6 Hz);
7.12 (2H, doublet of doublets, J=9 Hz, 5 Hz);
6.96 (2H, triplet, J=9 Hz);
6.32 (1H, doublet, J=3 Hz);
5.45 (1H, singlet);
2.99 (2H, broad doublet, J=3 Hz);
2.56 (2H, triplet, J=6 Hz);
2.39 (2H, triplet, J=8 Hz);
2.28-2.23 (2H, multiplet);
1.55-1.48 (2H, multiplet);
1.31-1.22 (10H, multiplet);
0.88 (3H, triplet, J=7 Hz).

Example 76

(±)-4-(1-Ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2667)

[0556]



[0557] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-1-ethyl-2-methylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl) isomer thereof (which is the compound of Example 77 below). The mixture was separated by chromatography

on a silica gel column using a 50:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 33%, R_f value = 0.20) as a pale yellow powder.

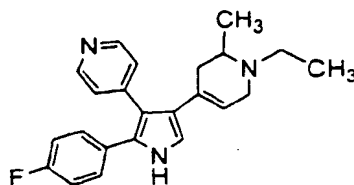
[0558] Melting point: 210 - 212°C (decomposition)

[0559] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
 8.31 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);
 7.13 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.97 (2H, triplet, J=9 Hz);
 6.32 (1H, singlet);
 5.27 (1H, singlet);
 3.09-3.00 (1H, multiplet);
 2.97-2.68 (1H, multiplet);
 2.34-2.73 (1H, multiplet);
 2.48-2.24 (3H, multiplet);
 2.22-2.11 (1H, multiplet);
 1.09 (3H, triplet, J=7 Hz);
 1.04 (3H, doublet, J=7 Hz).

Example 77

(±)-4-(1-Ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2709)



[0561] The title compound (yield 46%, R_f value = 0.15) was obtained as an orange powder during the chromatography performed in Example 76.

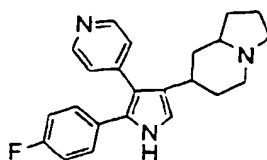
[0562] Melting point: 218 - 220°C (decomposition)

[0563] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
 8.32 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);
 7.12 (2H, doublet of doublets, J=9 Hz, 5 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.32 (1H, singlet);
 5.44 (1H, singlet);
 3.16-2.99 (2H, multiplet);
 2.90-2.79 (1H, multiplet);
 2.76-2.64 (1H, multiplet);
 2.51-2.32 (2H, multiplet);
 2.03-1.98 (1H, multiplet);
 1.08 (3H, triplet, J=7 Hz);
 1.01 (3H, doublet, J=6 Hz).

Example 78(±)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,7,8,8a-octahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2655)

[0564]



[0565] In a similar manner to that described in Example 11 above, (±)-2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 59 above) was reduced to give the title compound (yield 52%) as a white powder.

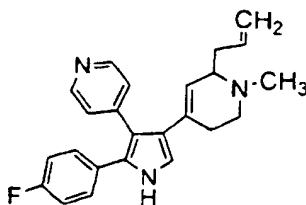
[0566] Melting point: 263 - 265°C

[0567] ¹H-Nuclear magnetic resonance spectrum (500 MHz, DMSO-d₆) δ ppm:

11.20 (1H, broad singlet);
 8.48 (2H, doublet, J=6 Hz);
 7.18-7.12 (4H, multiplet);
 7.09 (2H, triplet, J=9 Hz);
 6.76 (1H, doublet, J=2 Hz);
 2.98-2.92 (1H, multiplet);
 2.92-2.86 (1H, broad triplet, J=8 Hz);
 2.46 (1H, triplet of triplets, J=13Hz, 3 Hz);
 1.98-1.45 (9H, multiplet);
 1.30-1.14 (2H, multiplet).

Example 79(±)-4-(6-Allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2690)

[0568]



[0569] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-2-allyl-1-methylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (2-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl) isomer thereof (which is the compound of Example 80 below). The mixture was separated by chromatography on a silica gel column using a 100:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 20%, R_f value = 0.40) as a pale yellow powder.

[0570] Melting point : 218 - 220°C

[0571] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

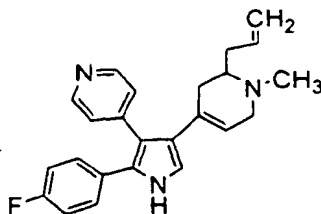
8.48 (2H, doublet, J=6 Hz);
 8.39 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);

7.11 (2H, doublet of doublets, J=9Hz, 5 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.33 (1H, doublet, J=3 Hz);
 5.66-5.58 (1H, multiplet);
 5.34 (1H, singlet);
 4.96 (1H, doublet, J=10 Hz);
 4.90 (1H, doublet, J=17 Hz);
 2.39-2.67 (1H, multiplet);
 2.71-2.67 (1H, multiplet);
 2.46-2.43 (2H, multiplet);
 2.37 (3H, singlet);
 2.35-2.32 (1H, multiplet);
 2.12-2.07 (2H, multiplet).

Example 80

(±)-4-(2-Allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2732)

[0572]



[0573] The title compound (yield 18%, Rf value = 0.35) was obtained as a pale yellow powder during the chromatography performed in Example 79 above.

[0574] Melting point : 178 - 180°C

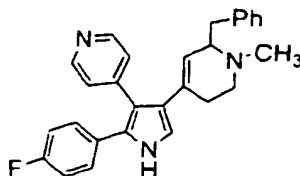
[0575] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
 8.38 (1H, broad singlet);
 7.17 (2H, doublet, J=6 Hz);
 7.12 (2H, doublet of doublets, J=9Hz, 5 Hz);
 6.96 (2H, triplet, J=9 Hz);
 6.32 (1H, doublet, J=3 Hz);
 5.77-5.69 (1H, multiplet);
 5.39 (1H, singlet);
 5.03 (1H, doublet, J=4 Hz);
 5.00 (1H, singlet);
 3.19-3.15 (1H, multiplet);
 2.99-2.95 (1H, multiplet);
 2.57-2.41 (1H, multiplet);
 2.36 (3H, singlet);
 2.36-2.31 (1H, multiplet);
 2.24-2.09 (3H, multiplet).

Example 81

(±)-4-(6-Benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2696)

[0576]



[0577] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-2-benzyl-1-methylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (2-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl) isomer thereof (which is the compound of Example 82 below). The mixture was separated by chromatography on a silica gel column using a 100:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 11%, Rf value = 0.38) as a pale yellow powder.

[0578] Melting point : 198 - 200°C

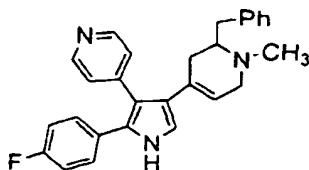
[0579] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.45 (2H, doublet, J=6 Hz);
 8.28 (1H, broad singlet);
 7.20-7.12 (3H, multiplet);
 7.09 (2H, doublet, J=6 Hz);
 7.10-7.07 (2H, multiplet);
 6.97-6.91 (4H, multiplet);
 6.76 (1H, doublet, J=3 Hz);
 5.24 (1H, singlet);
 3.13-3.10 (1H, multiplet);
 2.94-2.89 (2H, multiplet);
 2.51 (3H, singlet);
 2.49-2.35 (3H, multiplet);
 2.11-2.05 (1H, multiplet).

Example 82

(±)-4-(2-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2738)

[0580]



[0581] The title compound (yield 4%, Rf value = 0.33) was obtained as a pale yellow powder during the chromatography performed in Example 81 above.

[0582] Melting point: 198 - 199°C

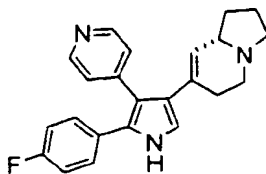
[0583] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.55 (1H, broad singlet);
 8.43 (2H, doublet, J=6 Hz);
 7.23 (2H, triplet, J=7 Hz);
 7.16 (2H, triplet, J=7 Hz);
 7.12 (2H, doublet, J=6 Hz);
 7.10-7.05 (4H, multiplet);
 6.94 (2H, triplet, J=9 Hz);
 6.72 (1H, doublet, J=3 Hz);
 5.40 (1H, singlet);
 3.20-3.17 (1H, multiplet);
 3.08-3.04 (2H, multiplet);
 2.31-2.78 (1H, multiplet);
 2.46 (3H, singlet);
 2.43-2.38 (1H, multiplet);
 2.10-2.00 (2H, multiplet).

Example 83

(-)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2653)

[0584]



[0585] (±)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (prepared as described in Example 60 above) was subjected to High Performance Liquid Chromatography using a chiral column and employing the conditions described below to give the title compound (retention time: 4.51 minutes, yield 94%) as a pale yellow powder.

[0586] Melting point: 204 - 205°C (Decomposition)

[0587] [α]_D²⁷ -50.67° (c=0.975, methanol)

<HPLC>

Column: CHIRALPAK AD (product of Daicel Chemical Industries, Ltd.)

Eluant: n-hexane : ethanol = 80:20

Flow rate: 1.0 ml/minute

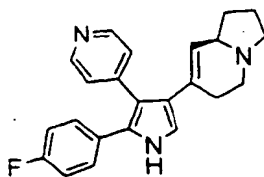
Temperature: 40°C

Detection: 254nm (UV).

Example 84

(+)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2653)

[0588]

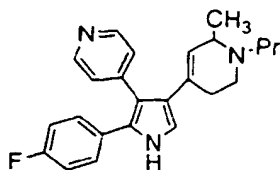


[0589] The title compound (retention time: 6.23 minutes, yield 95%) was obtained as a pale yellow powder during the High Performance Liquid Chromatography performed in Example 83 above.
 Melting Point: 207-210°C (Decomposition)
 $[\alpha]^{24}_D +46.59^\circ$ (c=0.920, methanol)

Example 85

(±)-2-(4-Fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2668)

[0590]



[0591] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-2-methyl-1-propylpiperidine-4-one as a starting material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl) isomer thereof (which is the compound of Example 86 below). The mixture was separated by chromatography on a silica gel column using a 100:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 6%, Rf value = 0.35) as a white powder.

[0592] Melting point : 214 - 215°C (decomposition)

[0593] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);

8.22 (1H, broad singlet);

7.17 (2H, doublet, J=6 Hz);

7.13 (2H, doublet of doublets, J=9Hz, 5 Hz);

6.97 (2H, triplet, J=9 Hz);

6.33 (1H, singlet);

5.27 (1H, singlet);

3.07-2.98 (1H, multiplet);

2.97-2.89 (1H, multiplet);

2.68-2.56 (1H, multiplet);

2.47-2.39 (1H, multiplet);

2.37-2.23 (2H, multiplet);

2.20-2.10 (1H, multiplet);

1.62-1.43 (2H, multiplet);

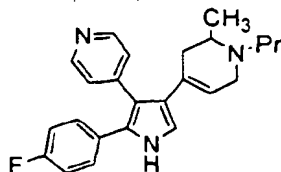
1.04 (3H, doublet, J=7 Hz);

0.90 (3H, triplet, J=7 Hz).

Example 86

(±)-2-(4-Fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2710)

[0594]



[0595] The title compound (yield 8%, Rf value = 0.32) was obtained as a white powder during the chromatography performed in Example 85 above.

[0596] Melting point : 210 - 218°C (decomposition)

[0597] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);

8.23 (1H, broad singlet);

7.17 (2H, doublet, J=6 Hz);

7.12 (2H, doublet of doublets J=9Hz, 5 Hz);

6.96 (2H, triplet, J=9 Hz);

6.82 (1H, singlet);

5.43 (1H, singlet);

3.16-3.01 (2H, multiplet);

2.89-2.82 (1H, multiplet);

2.58-2.50 (1H, multiplet);

2.42-2.30 (2H, multiplet);

2.00-1.92 (1H, multiplet);

1.57-1.45 (2H, multiplet);

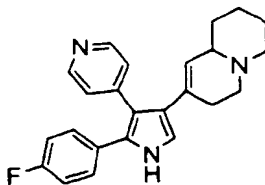
1.01 (3H, doublet, J=6 Hz);

0.90 (3H, triplet, J=7 Hz).

Example 87

(±)-2-(4-Fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2H-quinolizin-8-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2656)

[0598]



[0599] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using (±)-octahydro-2H-quinolizine-2-one as a starting material instead of 1-benzylpiperidine-4-one to afford a mixture of the title compound and the (1,3,4,6,9,9a-hexahydro-2H-quinolizin-8-yl) isomer thereof (which is the compound of Example 88 below). The mixture was separated by chromatography on a silica gel column using a 100:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 6%, Rf value = 0.35) as a pale yellow powder.

[0600] Melting point : 223 - 224°C

[0601] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);

8.30 (1H, broad singlet);

7.17 (2H, doublet, J=6 Hz);

7.12 (2H, doublet of doublets, J=9Hz, 5 Hz);

6.96 (2H, triplet, J=9 Hz);

6.33 (1H, doublet, J=3 Hz);

5.13 (1H, singlet);

2.38-2.61 (2H, multiplet);

2.67-2.56 (1H, multiplet);

2.45-2.43 (1H, multiplet);

2.40-2.35 (1H, multiplet);

2.17-2.07 (2H, multiplet);

1.76-1.73 (1H, multiplet);

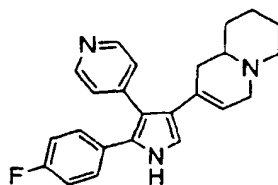
1.69-1.64 (2H, multiplet);

1.44-1.41 (1H, multiplet);

1.34-1.18 (2H, multiplet).

Example 88

(±)-2-(4-Fluorophenyl)-4-(1,3,4,6,9,9a-hexahydro-2H-quinolizin-8-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-2657)



[0603] The title compound (yield 19%, Rf value = 0.30) was obtained as a pale yellow powder during the chromatography performed in Example 87 above.

[0604] Melting point : 228 - 234°C (decomposition)

[0605] ¹H-Nuclear magnetic resonance spectrum (500 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);

8.28 (1H, broad singlet);

7.17 (2H, doublet, J=6 Hz);

7.12 (2H, doublet of doublets, J=9Hz, 5 Hz);

6.96 (2H, triplet, J=9 Hz);

6.31 (1H, doublet, J=3 Hz);

5.42 (1H, singlet);

3.30-3.25 (1H, multiplet);

2.98-2.95 (1H, multiplet);

2.77-2.73 (1H, multiplet);

2.14-1.99 (5H, multiplet);

1.74-1.65 (3H, multiplet);

1.31-1.21 (2H, multiplet).

Example 89

(±)-2-(4-Fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole
 (Compound No. 1-2906)

[0606]



[0607] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1-triisopropylsilyl-1H-pyrrole (prepared as described in Preparative Example 1 below) instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole, and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 90 below). The mixture was separated by chromatography on a silica gel column using a 200:20:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give the title compound (yield 15%, R_f value = 0.55) as a pale brown amorphous solid.

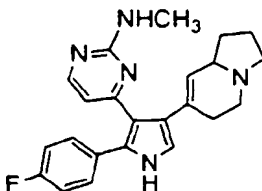
[0608] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CD₃OD) δ ppm:

8.08 (1H, doublet, J=5 Hz);
 7.30 (2H, doublet of doublets, J=9Hz, 5 Hz);
 7.02 (2H, triplet, J=9 Hz);
 6.32 (1H, singlet);
 6.49 (1H, doublet, J=5 Hz);
 5.57-5.51 (1H, multiplet);
 3.53-3.45 (1H, multiplet);
 3.21-3.11 (1H, multiplet);
 2.90-2.81 (1H, multiplet);
 2.30 (3H, singlet);
 2.47-2.28 (2H, multiplet);
 2.26-2.13 (2H, multiplet);
 2.07-1.96 (1H, multiplet);
 1.91-1.73 (2H, multiplet);
 1.49-1.37 (1H, multiplet).

Example 90

(±)-2-(4-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole
 (Compound No. 1-2905)

[0609]



[0610] The title compound (yield 12%, Rf value = 0.50) was obtained as a pale brown amorphous solid during the chromatography performed in Example 89 above.

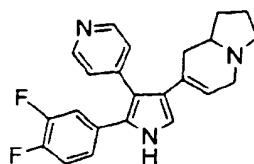
[0611] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CD₃OD) δ ppm:

8.09 (1H, doublet, J=5 Hz);
 7.30 (2H, doublet of doublets, J=9Hz, 5 Hz);
 7.02 (2H, triplet, J=9 Hz);
 6.33 (1H, singlet);
 6.46 (1H, doublet, J=5 Hz);
 5.48-5.42 (1H, multiplet);
 3.47-3.38 (1H, multiplet);
 3.03-2.75 (4H, multiplet);
 2.31 (3H, singlet);
 2.46-2.36 (1H, multiplet);
 2.29-2.20 (1H, multiplet);
 2.03-1.67 (2H, multiplet);
 1.33-1.75 (1H, multiplet);
 1.54-1.43 (1H, multiplet).

Example 91

(±)-2-(3,4-Difluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3260)

[0612]



[0613] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-(3,4-difluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole (prepared as described in Preparative Example 5 below), instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole, and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 92 below). The mixture was separated by chromatography on a silica gel column using a 100:10:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give the title compound (yield 29%, Rf value = 0.66) as a pale pink powder.

[0614] Melting point: 215 - 217°C (decomposition)

[0615] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

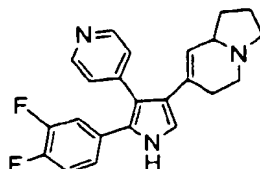
11.48 (1H, broad singlet);
 8.48 (2H, doublet, J=6 Hz);
 7.37-7.28 (1H, multiplet);
 7.21-7.10 (1H, multiplet);
 7.14 (2H, doublet, J=6 Hz);
 6.96 (1H, doublet, J=3 Hz);
 6.93-6.86 (1H, multiplet);
 5.27-5.20 (1H, multiplet);
 3.33-3.23 (1H, multiplet);
 3.05-2.97 (1H, multiplet);
 2.65-2.56 (1H, multiplet);
 2.30-2.22 (1H, multiplet);
 2.12-1.93 (3H, multiplet);
 1.91-1.81 (1H, multiplet);

1.75-1.57 (2H, multiplet);
1.33-1.20 (1H, multiplet).

Example 92

(±)-2-(3,4-Difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3259)

[0616]



[0617] The title compound (yield 23%, Rf value = 0.31) was obtained as a pale pink powder during the chromatography performed in Example 91 above.

[0618] Melting point: 187 - 190°C (decomposition)

[0619] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

11.42 (1H, broad singlet);

8.47 (2H, doublet, J=6 Hz);

7.39-7.29 (1H, multiplet);

7.21-7.10 (1H, multiplet);

7.14 (2H, doublet, J=6 Hz);

6.96 (1H, doublet, J=3 Hz);

6.94-6.87 (1H, multiplet);

5.18-5.13 (1H, multiplet);

3.10-3.02 (1H, multiplet);

2.89-2.82 (1H, multiplet);

2.79-2.71 (1H, multiplet);

2.63-2.45 (2H, multiplet);

2.30-2.20 (1H, multiplet);

2.06-1.97 (1H, multiplet);

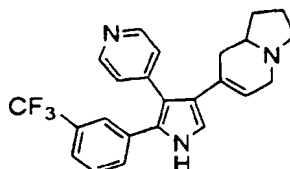
1.78-1.53 (3H, multiplet);

1.22-1.10 (1H, multiplet).

Example 93

(±)-4-(1,2,3,5,6,8a-Hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1H-pyrrole (Compound No. 1-3262)

[0620]



[0621] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1-triisopropylsilyl-1H-

pyrrole (prepared as described in Preparative Example 6 below), instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1*H*-pyrrole, and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 94 below). The mixture was separated by chromatography on a silica gel column using a 100:10:1 by volume mixture of ethyl acetate, methanol and isopropylamine as the eluant to give the title compound (yield 26%, R_f value = 0.65) as a pale brown powder.

[0622] Melting point: 220 - 223°C (decomposition)

[0623] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

11.62 (1H, broad singlet);

8.49 (2H, doublet, J=6 Hz);

7.54-7.35 (4H, multiplet);

7.16 (2H, doublet, J=6 Hz);

7.02 (1H, doublet, J=3 Hz);

5.29-5.21 (1H, multiplet);

3.35-3.24 (1H, multiplet);

3.06-2.98 (1H, multiplet);

2.65-2.56 (1H, multiplet);

2.34-2.24 (1H, multiplet);

2.13-1.94 (3H, multiplet);

1.92-1.82 (1H, multiplet);

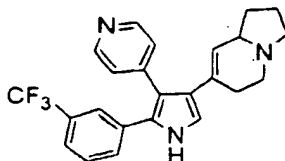
1.75-1.58 (2H, multiplet);

1.34-1.21 (1H, multiplet).

Example 94

(±)-4-(1,2,3,5,6,8a-Hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1*H*-pyrrole (Compound No. 1-3261)

[0624]



[0625] The title compound (yield 17%, R_f value = 0.31) was obtained as a pale brown powder during the chromatography performed in Example 93 above.

[0626] Melting point : 189 - 191°C (decomposition)

[0627] ¹H-Nuclear magnetic resonance spectrum (400 MHz, DMSO-d₆) δ ppm:

11.63 (1H, broad singlet);

8.48 (2H, doublet, J=6 Hz);

7.55-7.37 (4H, multiplet);

7.15 (2H, doublet, J=6 Hz);

7.02 (1H, doublet, J=3 Hz);

5.19-5.14 (1H, multiplet);

3.12-3.03 (1H, multiplet);

2.91-2.82 (1H, multiplet);

2.30-2.71 (1H, multiplet);

2.64-2.45 (2H, multiplet);

2.33-2.21 (1H, multiplet);

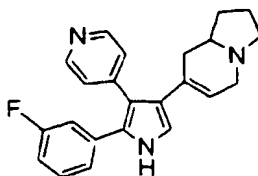
2.09-1.98 (1H, multiplet);

1.78-1.53 (3H, multiplet);

1.21-1.10 (1H, multiplet).

Example 95(±)-2-(3-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3256)

[0628]



[0629] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-(3-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole (prepared as described in Preparative Example 4 below), instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole, and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 96 below). The mixture was separated by chromatography on a silica gel column using a 30:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 17%, Rf value = 0.25) as a pale pink powder.

[0630] Melting point : 199 - 203°C (decomposition)

[0631] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.49 (2H, doublet, J=6 Hz);

8.30 (1H, broad singlet);

7.23-7.18 (3H, multiplet);

6.96-6.82 (4H, multiplet);

5.50-5.47 (1H, multiplet);

3.53-2.96 (1H, multiplet);

3.20 (1H, triplet, J=9 Hz);

2.79 (1H, doublet, J=17 Hz);

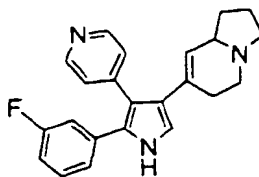
2.32-2.08 (4H, multiplet);

2.00-1.67 (3H, multiplet);

1.48-1.33 (1H, multiplet).

Example 96(±)-2-(3-Fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3255)

[0632]



[0633] The title compound (yield 12%, Rf value = 0.10) was obtained as a pale brown powder during the chromatography performed in Example 95 above.

[0634] Melting point : 178 - 181°C (decomposition)

[0635] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

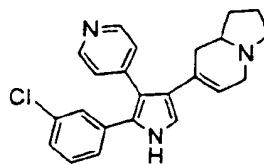
8.50 (2H, doublet, J=6 Hz);

8.34 (1H, broad singlet);

7.26-7.18 (3H, multiplet);
 6.95-6.84 (4H, multiplet);
 5.42 (1H, singlet);
 3.23-3.17 (1H, multiplet);
 5 3.02-2.82 (2H, multiplet);
 2.78-2.61 (2H, multiplet);
 2.42-2.30 (1H, multiplet);
 2.16-2.04 (1H, multiplet);
 1.95-1.68 (3H, multiplet);
 10 1.44-1.32 (1H, multiplet).

Example 97

(±)-2-(3-Chlorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3258)

[0636]

[0637] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-(3-chlorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole (prepared as described in Preparative Example 3 below), instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 98 below). The mixture was separated by chromatography on a silica gel column using a 30:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 18%, Rf value = 0.25) as a pale pink powder.

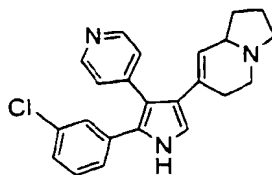
[0638] Melting point : 197 - 201°C (decomposition)

[0639] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.49 (2H, doublet, J=6 Hz);
 8.32 (1H, broad singlet);
 7.25-7.13 (5H, multiplet);
 6.96 (1H, doublet, J=7 Hz);
 6.35 (1H, doublet, J=3 Hz);
 5.50-5.47 (1H, multiplet);
 3.54-3.45 (1H, multiplet);
 3.20 (1H, triplet, J=9 Hz);
 2.79 (1H, doublet, J=16 Hz);
 2.31-2.07 (4H, multiplet);
 2.00-1.67 (3H, multiplet);
 1.48-1.33 (1H, multiplet).

Example 98(±)-2-(3-Chlorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3257)

[0640]



[0641] The title compound (yield 16%, Rf value = 0.10) was obtained as a pale brown powder during the chromatography performed in Example 97 above.

[0642] Melting point : 193 - 195°C (decomposition)

[0643] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.50 (2H, doublet, J=6 Hz);

8.42 (1H, broad singlet);

7.25-7.13 (5H, multiplet);

6.96 (1H, doublet, J=7 Hz);

6.35 (1H, doublet, J=3 Hz);

5.42 (1H, singlet);

3.26-3.17 (1H, multiplet);

3.02-2.82 (2H, multiplet);

2.78-2.60 (2H, multiplet);

2.42-2.30 (1H, multiplet);

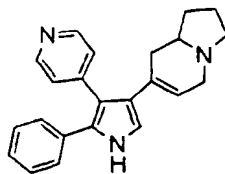
2.16-2.03 (1H, multiplet);

1.95-1.65 (3H, multiplet);

1.45-1.32 (1H, multiplet).

Example 99(±)-4-(1,2,3,5,8,8a-Hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3254)

[0644]



[0645] In a similar manner to the procedures described in Examples 9(i) and 9(iii) above, coupling, dehydration and desilylation reactions were carried out using 4-bromo-2-phenyl-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole (prepared as described in Preparative Example 2 below), instead of 4-bromo-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole, and (±)-1,2,3,5,6,7,8,8a-octahydroindolizine-7-one, instead of 1-benzylpiperidine-4-one, as starting materials to afford a mixture of the title compound and the (1,2,3,5,6,8a-hexahydroindolizin-7-yl) isomer thereof (which is the compound of Example 100 below). The mixture was separated by chromatography on a silica gel column using a 30:1 by volume mixture of ethyl acetate and isopropylamine as the eluant to give the title compound (yield 14%, Rf value = 0.24) as a pale brown powder.

[0646] Melting point: 201 - 204°C (decomposition)

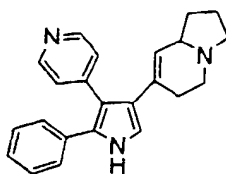
[0647] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.46 (2H, doublet, J=6 Hz);
 8.28 (1H, broad singlet);
 7.30-7.13 (7H, multiplet);
 6.34 (1H, doublet, J=3 Hz);
 5.50-5.47 (1H, multiplet);
 3.54-3.45 (1H, multiplet);
 3.50 (1H, triplet, J=9 Hz);
 2.79 (1H, doublet, J=17 Hz);
 2.32-2.07 (4H, multiplet);
 2.00-1.67 (3H, multiplet);
 1.48-1.33 (1H, multiplet).

Example 100

(±)-4-(1,2,3,5,6,8a-Hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1H-pyrrole (Compound No. 1-3253)

[0648]



[0649] The title compound (yield 10%, R_f value = 0.11) was obtained as a pale brown powder during the chromatography performed in Example 99 above.

[0650] Melting point : 180 - 183°C (decomposition)

[0651] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.47 (2H, doublet, J=6 Hz);
 8.34 (1H, broad singlet);
 7.30-7.13 (7H, multiplet);
 6.34 (1H, doublet, J=3 Hz);
 5.44 (1H, singlet);
 3.27-3.18 (1H, multiplet);
 3.02-2.82 (2H, multiplet);
 2.78-2.60 (2H, multiplet);
 2.43-2.30 (1H, multiplet);
 2.17-2.06 (1H, multiplet);
 1.95-1.65 (3H, multiplet);
 1.45-1.32 (1H, multiplet).

Preparative Examples

Preparative Example 1

4-Bromo-2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1-triisopropylsilyl-1H-pyrrole

1(i) 2-(4-Fluorophenyl)-3-(2-methylthiopyrimidin-4-yl)-1H-pyrrole

[0652] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction) and Example 7(i) above (decarboxylation), reactions were carried out using ethyl 3-(2-methylthiopyrimidin-4-yl)acrylate instead of ethyl 3-(4-pyridyl)acrylate as a starting material to afford the title compound (yield 37%) as a white powder.

[0653] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.41 (1H, broad singlet);
 8.24 (1H, doublet, J=5 Hz);

7.42 (1H, doublet of doublets, J=9Hz, 5 Hz);
 7.10 (2H, triplet, J=9 Hz);
 6.38 (1H, triplet, J=3 Hz);
 6.34 (1H, doublet, J=5 Hz);
 5 6.31 (1H, triplet, J=3 Hz);
 2.33 (3H, singlet).

1(ii) 2-(4-Fluorophenyl)-3-(2-methanesulfonylpyrimidin-4-yl)-1H-pyrrole

10 **[0654]** In a similar manner to that described in Example 42(ii) above, 2-(4-fluorophenyl)-3-(2-methylthiopyrimidin-4-yl)-1H-pyrrole [which was prepared as described in step 1(i) above] was oxidized to afford the title compound (yield 81%) as a pale yellow powder.

[0655] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃-CD₃OD) δ ppm:

8.58 (1H, doublet, J=6 Hz);
 15 7.50 (2H, doublet of doublets, J=9Hz, 6 Hz);
 7.47 (1H, doublet, J=6 Hz);
 7.16 (2H, triplet, J=9 Hz);
 6.90 (1H, doublet, J=3 Hz);
 6.34 (1H, doublet, J=3 Hz);
 20 3.05 (3H, singlet).

1[(iii) 2-(4-Fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole

25 **[0656]** In a similar manner to that described in Example 42(iii) above, 2-(4-fluorophenyl)-3-(2-methanesulfonylpyrimidin-4-yl)-1H-pyrrole [prepared as described in step 1(ii) above] was reacted with methylamine to afford the title compound (quantitative yield) as a pale brown amorphous solid.

[0657] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.37 (1H, broad singlet);
 8.08 (1H, doublet, J=5 Hz);
 30 7.44 (2H, doublet of doublets, J=9Hz, 6 Hz);
 7.08 (2H, triplet, J=9 Hz);
 6.35 (1H, triplet, J=3 Hz);
 6.75 (1H, triplet, J=3 Hz);
 6.47 (1H, doublet, J=5 Hz);
 35 4.90 (1H, broad quartet, J=5 Hz);
 2.37 (3H, doublet, J=5 Hz).

1(iv) 2-(4-Fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1-triisopropylsilyl-1H-pyrrole

40 **[0658]** In a similar manner to that described in Example 7(ii) above, 2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole [prepared as described in step 1(iii) above] was reacted with triisopropylsilyl triflate, to afford the title compound (yield 82%) as a white powder.

[0659] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

7.92 (1H, doublet, J=5 Hz);
 45 7.36 (2H, doublet of doublets, J=9Hz, 6 Hz);
 7.06 (2H, triplet, J=9 Hz);
 6.90 (1H, doublet, J=3 Hz);
 6.89 (1H, doublet, J=3 Hz);
 5.93 (1H, doublet, J=5 Hz);
 50 4.30 (1H, broad quartet, J=5 Hz);
 2.33 (3H, doublet, J=5 Hz);
 1.15-0.92 (21H, multiplet).

1(v) 3-[2-(N-t-Butoxycarbonyl-N-methylamino)pyrimidin-4-yl]-2-(4-fluorophenyl)-1-triisopropylsilyl-1H-pyrrole

55 **[0660]** In a similar manner to that described in Example 4(i) above, 2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 1(iv) above] was reacted with di-t-butyl dicarbonate to afford the title compound (yield 90%) as colorless oil.

[0661] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

- 8.24 (1H, doublet, J=5 Hz);
 7.36 (2H, doublet of doublets, J=9Hz, 5 Hz);
 7.09 (2H, triplet, J=9 Hz);
 5 6.95 (1H, doublet, J=3 Hz);
 6.90 (1H, doublet, J=3 Hz);
 6.28 (1H, doublet, J=5 Hz);
 3.20 (3H, singlet);
 1.51 (9H, singlet);
 10 1.15-0.94 (21H, multiplet).

1(vi) 4-Bromo-3-[2-(N-t-butoxycarbonyl-N-methylamino)pyrimidin-4-yl]-2-(4-fluorophenyl)-1-triisopropylsilyl-1H-pyrrole

- 15 **[0662]** In a similar manner to that described in Example 7(iii) above, 3-[2-(N-t-butoxycarbonyl-N-methylamino)pyrimidin-4-yl]-2-(4-fluorophenyl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step 1(v) above] was brominated to afford the title compound (yield 79%) as a pale brown amorphous solid.

[0663] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

- 8.42 (1H, doublet, J=5 Hz);
 20 7.28 (2H, doublet of doublets, J=9Hz, 5 Hz);
 7.00 (2H, triplet, J=9 Hz);
 6.93 (1H, singlet);
 6.78 (1H, doublet, J=5 Hz);
 3.13 (3H, singlet);
 25 1.49 (9H, singlet);
 1.15-0.93 (21H, multiplet).

1(vii) 4-Bromo-2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole

- 30 **[0664]** 472 mg (0.782 mmol) of 4-bromo-3-[2-(N-t-butoxycarbonyl-N-methylamino)pyrimidin-4-yl]-2-(4-fluorophenyl)-1-triisopropylsilyl-1H-pyrrole [prepared as described in step (vi) above] were dissolved in 10 of ml tetrahydrofuran, 0.98 ml (3.91 mmol) of a 4N solution of hydrogen chloride in dioxane were added to the solution, and the resulting mixture was stirred for 3 hours at 50 °C. At the end of this time, water was added to the reaction mixture and then it was
 35 neutralized with a saturated aqueous solution of sodium hydrogencarbonate. This was extracted with ethyl acetate, and the organic extract was concentrated under reduced pressure. The resulting residue was purified by chromatography on a silica gel column using a 7:3 by volume mixture of hexane and ethyl acetate as the eluant to afford 202 mg (yield 74%) of the title compound as a pale brown amorphous solid.

[0665] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

- 8.66 (1H, broad singlet);
 40 8.21 (1H, doublet, J=5 Hz);
 7.26 (2H, doublet of doublets, J=9Hz, 5 Hz);
 7.01 (2H, triplet, J=9 Hz);
 6.38 (1H, doublet, J=3 Hz);
 6.61 (1H, doublet, J=5 Hz);
 45 4.96 (1H, broad quartet, J=5 Hz);
 2.86 (3H, doublet, J=5 Hz).

1 (viii) 4-Bromo-2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1-triisopropylsilyl-1H-pyrrole

- 50 **[0666]** In a similar manner to the procedures described in Example 7(ii) above, 4-bromo-2-(4-fluorophenyl)-3-(2-methylaminopyrimidin-4-yl)-1H-pyrrole [prepared as described in step 1(vii) above] was reacted with triisopropylsilyl triflate to afford the title compound (yield 67%) as a pale brown amorphous solid.

[0667] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

- 8.04 (1H, doublet, J=5 Hz);
 55 7.30 (2H, doublet of doublets, J=9Hz, 6 Hz);
 6.98 (2H, triplet, J=9 Hz);
 6.91 (1H, singlet);
 6.32 (1H, broad singlet);

4.35 (1H, broad quartet, J=4 Hz);
 2.32 (3H, doublet, J=4 Hz);
 1.15-0.92 (21H, multiplet).

Preparative Example 2

4-Bromo-2-phenyl-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0668] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction), Example 7(i) above (decarboxylation), Example 7(ii) above (reaction with triisopropylsilyl triflate) and Example 7(iii) above (bromination), reactions were carried out using α -(p-toluenesulfonyl)benzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

[0669] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.35 (2H, doublet, J=6 Hz);
 7.36-7.30 (1H, multiplet);
 7.28-7.21 (4H, multiplet);
 7.03 (2H, doublet, J=6 Hz);
 6.94 (1H, singlet);
 1.15-1.02 (21H, multiplet).

Preparative Example 3

4-Bromo-2-(3-chlorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0670] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction), Example 7(i) above (decarboxylation), Example 7(ii) above (reaction with triisopropylsilyl triflate) and Example 7(iii) above (bromination), reactions were carried out using α -(p-toluenesulfonyl)-3-chlorobenzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

[0671] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.40 (2H, doublet, J=6 Hz);
 7.33-7.29 (2H, multiplet);
 7.18 (1H, triplet, J=8 Hz);
 7.09 (1H, doublet, J=8 Hz);
 7.03 (2H, doublet, J=6 Hz);
 6.95 (1H, singlet);
 1.12-0.99 (21H, multiplet).

Preparative Example 4

4-Bromo-2-(3-fluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0672] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction), Example 7(i) above (decarboxylation), Example 7(ii) above (reaction with triisopropylsilyl triflate) and Example 7(iii) above (bromination), reactions were carried out using α -(p-toluenesulfonyl)-3-fluorobenzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

[0673] ¹H-Nuclear magnetic resonance spectrum (400 MHz, CDCl₃) δ ppm:

8.39 (2H, doublet, J=6 Hz);
 7.22 (1H, doublet of doublets, J=8Hz, 6 Hz);
 7.07-7.01 (4H, multiplet);
 6.98-6.94 (2H, multiplet);
 1.14-0.98 (21H, multiplet).

Preparative Example 5

4-Bromo-2-(3,4-difluorophenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0674] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction), Example 7(i) above (decarboxylation), Example 7(ii) above (reaction with triisopropylsilyl triflate) and Example 7(iii) above (bromination), reactions were carried out using α -(p-toluenesulfonyl)-3,4-difluorobenzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

mination), reactions were carried out using α -(p-toluenesulfonyl)-3,4-difluorobenzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

[0675] ^1H -Nuclear magnetic resonance spectrum (400 MHz, CDCl_3) δ ppm:

8.42 (2H, doublet, $J=6$ Hz);

7.12-6.93 (3H, multiplet);

7.02 (2H, doublet, $J=6$ Hz);

6.94 (1H, singlet);

1.18-0.94 (21H, multiplet).

10 Preparative Example 6

4-Bromo-2-(3-trifluoromethylphenyl)-3-(pyridin-4-yl)-1-triisopropylsilyl-1H-pyrrole

[0676] In a similar manner to the procedures described in Example 1(i) above (pyrrole ring forming reaction), Example 7(i) above (decarboxylation), Example 7(ii) above (reaction with triisopropylsilyl triflate) and Example 7(iii) above (bromination), reactions were carried out using α -(p-toluenesulfonyl)-3-trifluoromethylbenzylisocyanide, instead of α -(p-toluenesulfonyl)-4-fluorobenzylisocyanide, as a starting material to afford the title compound as a pale purple powder.

[0677] ^1H -Nuclear magnetic resonance spectrum (400 MHz, CDCl_3) δ ppm:

8.39 (2H, doublet, $J=6$ Hz);

7.64-7.57 (2H, multiplet);

7.40-7.33 (2H, multiplet);

7.00 (2H, doublet, $J=6$ Hz);

6.98 (1H, singlet);

1.15-0.94 (21H, multiplet).

25 Formulation Examples

[0678] A pharmaceutical preparation containing a compound of the present invention having the above formula (I), or a pharmacologically acceptable salt, ester or other derivative thereof as its active ingredient can be produced according to, for example, the following methods.

Formulation Example 1

Powder

[0679] 5 g of the compound of Example 2, 895 g of lactose and 100 g of corn starch were mixed in a blender to provide the desired powder.

Formulation Example 2

Granules

[0680] 5 g of the compound of Example 5, 865 g of lactose and 100 g of low-substituted hydroxypropylcellulose were mixed, 300 g of a 10% aqueous hydroxypropyl cellulose solution were added to the resulting mixture, and this was then kneaded. The product thus obtained was then granulated using an extrusion granulating machine and dried to provide the desired granules.

Preparation Example 3

Capsules

[0681] 5 g of the compound of Example 6, 115 g of lactose, 58 g of corn starch and 2 g of magnesium stearate were mixed using a V-shaped mixer, no. 3 capsules were chosen and then each of said no. 3 capsules was filled with 180 mg of the resulting mixture to provide the desired capsules.

Preparation Example 4Tablets

[0682] 5 g of the compound of Example 83, 90 g of lactose, 34 g of corn starch, 20 g of crystalline cellulose and 1 g of magnesium stearate were mixed in a blender, and the resulting mixture was then formed into tablets with a tablet machine to provide the desired tablets.

Test Examples

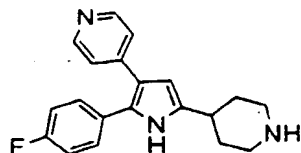
[0683] The biological activity of the compounds of the present invention is illustrated by the following Test Examples.

Test Example 1Inhibition of the Production of the Cytokines IL-1 β and TNF α *in vitro* in Human Whole Blood

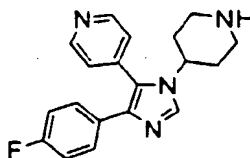
[0684] This test was performed according to the method of Hartman, et al. [D.A. Hartman, S.J. Ochalski and R.P. Carlson; The effects of anti-inflammatory and antiallergic drugs on cytokine release after stimulation of human whole blood by lipopolysaccharide and zymosan A: *Inflamm. Res.*, 44, 269 (1995)].

[0685] Peripheral blood samples were collected in the presence of heparin from healthy adult volunteers. 1000 μ l of whole blood were added to an Eppendorf tube to which 2 μ l of a dimethyl sulfoxide solution of the test compound had been added in advance, after which 10 μ l of lipopolysaccharide (E. coli 026: B6 origin, Difco) were added as a stimulant (final concentration of said lipopolysaccharide: 10 μ g/ml). This was mixed well and then incubated for 6 hours at 37°C in the presence of 5% CO₂. At the end of the incubation, the mixture was cooled to 4°C to stop the reaction, followed immediately by centrifuging for 5 minutes at 14,000 rpm to separate and collect the supernatant plasma. The IL-1 β and TNF α produced and released into the plasma were measured using a commercially available enzyme immunoassay (ELISA) kit [Cayman (IL-1 β) and Genzyme (TNF α)]. The procedure was also repeated in the absence of test compound. The inhibitory effect [IC₅₀ (μ M)] on the production of IL-1 β and TNF α was determined from the amounts of the cytokines produced in the presence and absence of the test compound.

[0686] The results for the inhibition of the production *in vitro* of TNF α are shown in Table 3 below. In this Table, Compounds A and B are as follows:



Compound A
(Compound of Example 4
of WO97/5877)



Compound B
(Compound of Example 23
of WO96/21452)

Table 3

Example No.	Inhibitory Effect on TNF α Production [IC ₅₀ (μ M)]
Compound A	1.90
Compound B	1.73
45	0.14
58	0.089
60	0.047
83	0.026

Table 3 (continued)

Example No.	Inhibitory Effect on TNF α Production [IC ₅₀ (μ M)]
85	0.44
86	0.29
87	0.31
92	0.045
94	0.031

[0687] Table 3 above demonstrates excellent inhibitory activity against the production of TNF α *in vitro* for the compounds of the present invention.

[0688] The compounds of the present invention were also found to show excellent inhibitory activity against the production of IL-1 β *in vitro*

Test Example 2

Inhibition of the Production of TNF α *in vivo*

[0689] This test was performed according to the method of Ochalski, et al. [S.J. Ochalski, D.A. Hartman, M.T. Belfast, T.L. Walter, K.B. Glaser and R.P. Carlson; Inhibition of endotoxin-induced hypothermia and serum TNF- α levels in CD-1 mice by various pharmacological agents: Agents Actions 39, C52-C54 (1993)].

[0690] The production of TNF α was induced in mice by the intravenous injection of lipopolysaccharide (E. coli O26: B6 origin, Difco) which was prepared to a concentration of 0.045 mg/ml using physiological saline. The saline preparation of lipopolysaccharide was administered at the rate of 10 ml/1 kg of body weight into the caudal vein of Balb/c mice (males, age 5-7 weeks, body weight: approx. 22 g, Japan Charles River) which had been fasted overnight starting on the day before the experiment. One hour after administration, the mice were laparotomized under ether anaesthesia and blood was collected from the abdominal vena cava. Blood collection was performed using a 1 ml volume disposable syringe equipped with a 23G needle which had been moistened with heparin on the inside wall. Following blood collection, the blood was immediately transferred to a 1.5 ml volume Eppendorf tube and centrifuged at 4°C and 14,000 rpm to separate the plasma. This plasma was then stored at -20°C until measurement of TNF α . The measurement of the amount of TNF α was performed with a commercially available enzyme immunoassay (ELISA) kit (Mouse TNF α ELISA KIT, Genzyme).

[0691] To determine the inhibitory activity of the test compounds, each test compound was suspended in a 0.5% tragacanth solution and then administered orally to the Balb/c mice at the rate of 10 ml/1 kg of body weight 30 minutes before injection of lipopolysaccharide. The level of TNF α production was then determined as described above. In the control group, 0.5% tragacanth solution was administered at the rate of 10 ml/1 kg of body weight to the test mice instead of the solutions of the test compounds. A minimum of 3 dose levels of the test compound was administered to groups of 5 test mice for each test compound. The inhibitory rate relative to the control group was calculated for each dose level. From the inhibitory rates and the dosages, ID₅₀ values were calculated by the least squares method.

[0692] The compounds of the present invention were found to show excellent inhibitory activity against the production of TNF α *in vivo*.

Test Example 3

Inhibition of the Production of IL-1 β *in vivo*

[0693] This test was performed according to the method of Griffiths, et al. [Richard J. Griffiths, Ethan J. Stam, James T. Downs and Ivan G. Otterness; ATP Induces the Release of IL-1 from LPS-Primed Cells *In Vivo*: J. Immunol., 154, 2821-2828 (1995)].

[0694] The production of IL-1 was induced in mice by the intraperitoneal injection of lipopolysaccharide followed by the intraperitoneal injection of adenosine triphosphate (ATP). This was achieved by first administering a solution of lipopolysaccharide (E. coli O26: B6 origin, Difco), which had been prepared to a concentration of 0.0045 mg/ml using physiological saline, at the rate of 10 ml of said saline solution/1 kg of body weight into the peritoneal cavity of Balb/c mice (males, age 5-7 weeks, body weight: approx. 22 g, Japan Charles River) which had been fasted overnight starting on the day before the experiment. Two hours later, 0.5 ml of ATP, which had been prepared to a concentration of 6.03

mg/ml using physiological saline, were administered into the peritoneal cavity. 0.5 hours after the administration of ATP, the mice were sacrificed by suffocation using dry ice followed immediately by intraperitoneal injection of 3 ml of washing phosphate buffer solution [containing heparin (10 U/ml), p-toluenesulfonyl fluoride (0.25 mM), leupepsin (1 µg/ml), pepstatin (1 µg/ml) and EDTA (1 mM)] to wash the peritoneal cavity. A 1 ml volume disposable syringe equipped with a 21G needle was then used to recover the washing liquid. After the recovery, the washing liquid from the peritoneal cavity was immediately transferred to a 1.5 ml volume Eppendorf tube and centrifuged at 4°C and 7,500 rpm to separate the supernatant. This supernatant was then stored at -20°C until measurement of IL-1β.

[0695] The measurement of the amount of IL-1β was performed with an enzyme immunoassay (ELISA) kit (Mouse IL-1β ELISA KIT, Genzyme).

[0696] To determine the inhibitory activity of the test compounds, each test compound was suspended in a 0.5% tragacanth solution and then administered orally to the Balb/c mice at the rate of 10 ml/kg of body weight 30 minutes before injection of lipopolysaccharide. The level of TNFα production was then determined as described above. In the control group 0.5% tragacanth solution was administered to the test mice at the rate of 10 ml/kg of body weight instead of the solutions of the test compounds. A minimum of 3 dose levels of the test compound was administered to groups of 5 test mice for each test compound. The mean inhibitory rate relative to the control group was calculated for each dose level.

[0697] In this test the compounds of the present invention demonstrated an excellent inhibitory effect against the production of IL-1β *in vivo*.

Test Example 4

Activity in Preventing the Development of Adjuvant-Induced Arthritis *in vivo*

[0698] The test was performed according to the method described by Winder et al. (Arthritis Rheum., 12, 472-482, 1969).

[0699] Heat-killed dried Mycobacterium butyricum (Difco Laboratories, Lot 679123) was ground on an agate mortar, and was then suspended in dry-sterilised liquid paraffin (first grade, Wako Pure Chemical Industries, Ltd.) to make a 2 ml suspension. The resulting suspension was then sonicated and used as an adjuvant. Arthritis was induced by the intradermal injection of the adjuvant (100 µg of heat killed dried bacterium/0.05 ml of paraffin/paw) into the heel of the right hind paw of a Lewis rat (male, age 8 weeks, Japan Charles River). The test compounds, which had been suspended in 0.5% sodium carboxymethyl cellulose solution (CMC, Daiichi Pure Chemicals, Co., Ltd.), were administered orally once a day from the day of injection of the adjuvant (day 0) to day 20.

[0700] The volumes of the injected (right) and non-injected (left) hind paws were measured on days 3, 5, 7, 10, 13, 15, 18 and 21 using a Plethysmometer™ (Ugo Basile), the hind paws being soaked from the toe to the hairline in the bath of the Plethysmometer™. The volumes of the swollen feet (adjuvant-injected right hind foot volume - non-injected left hind foot volume) were calculated. The percent inhibition of swelling of the injected foot of the treated animals as compared to that of the control animals on day 21 was calculated as follows.

$$\text{Inhibition (\%)} = \{1 - (\text{swollen foot volume of compound-treated animals}) / (\text{swollen foot volume of control animals})\} \times 100$$

[0701] A linear regression curve was obtained from the percent inhibition and the logarithmic value of the dosage by the least squares method. ID₅₀ values were calculated using this curve.

[0702] In this test, the compounds of the present invention showed excellent activity in preventing the development of adjuvant-induced arthritis.

Test Example 5

Activity in Preventing the Development of Arthritis Induced by Anti-Collagen Antibody *in vivo*

[0703] In this test, an anti-collagen antibody-induced mouse arthritis model was employed.

[0704] 0.5 ml (2 mg of antibody) of an anti-collagen antibody solution (4 mg/ml, Arthritogenic mAb Cocktail: product of Immuno-Biological Laboratories Co., Ltd) were injected into the caudal vein of Balb/c mice (males, age 5-6 weeks old, Japan Charles River). Three days after injection, 0.1 ml [0.05 mg of lipopolysaccharide] of a lipopolysaccharide solution (0.5 mg/ml, Arthritogenic mAb Cocktail: product of Immuno-Biological Laboratories Co., Ltd) were administered to the mice by intraperitoneal injection.

[0705] The test compounds, which had been suspended in 0.5% tragacanth were administered orally to the test animals at the rate of 10 ml/1 kg of body weight once per day for 7 days from the day when the anti-collagen antibody was administered. To the mice of the control group, 0.5% tragacanth solution was administered at the rate of 10 ml/kg of body weight once per day for 7 days from the day when the anti-collagen antibody was administered, instead of solutions of the test compounds.

[0706] After the administration of the test compounds (or 0.5% tragacanth solution), the degree of edema in the 4 paws of each test mouse was scored according to the following basis:

- 0: normal (edema is not observed);
- 1: edema is observed in one of the five toes;
- 2: edema is observed in two or more of the five toes;
- 3: the whole of the paw is swollen.

[0707] The degree of arthritis in the test mouse was evaluated by the total of the edema scores in the 4 paws. The rate of suppression was calculated from the degrees of arthritis of the control animals and of the animals treated with the test compounds. From the rates of suppression and the dosages, ID₅₀ values were calculated by the least squares method.

[0708] In this test, the compounds of the present invention showed excellent activity in preventing the development of arthritis induced by anti-collagen antibody.

Test Example 6

Activity in Treating Arthritis Induced by Anti-Collagen Antibody *in vivo*

[0709] In this test, an anti-collagen antibody-induced mouse arthritis model was employed.

[0710] 0.5 ml (2 mg of antibody) of an anti-collagen antibody solution (4 mg/ml, Arthritisogenic mAb Cocktail: product of Immuno-Biological Laboratories Co., Ltd) were injected into the caudal vein of Balb/c mice (males, age 5-6 weeks old, Japan Charles River). Three days after injection, 0.1 ml [0.05 mg of lipopolysaccharide] of a lipopolysaccharide solution (0.5 mg/ml, Arthritisogenic mAb Cocktail: product of Immuno-Biological Laboratories Co., Ltd) were administered to the mice by intraperitoneal injection.

[0711] 7 days after the administration of the anti-collagen antibody solution, the degree of edema in the 4 paws of each test mouse was scored according to the basis as shown in Test Example 5 above.

[0712] Those mice in which edema in both the hind paws had been scored as "3" were selected. Test compounds, which had been suspended in 0.5% tragacanth solution, were administered orally to the selected mice at the rate of 10 ml/kg of body weight once per day for 3 days. To the mice of the control group, 0.5% tragacanth solution was administered at the rate of 10 ml/kg of body weight once per day for 3 days instead of solutions of the test compounds.

[0713] After the administration of the test compounds (or 0.5% tragacanth solution), the degree of arthritis in each test mouse was evaluated in the same manner as described in Test Example 5. The rates of treatment of arthritis induced by anti-collagen antibody were calculated from the degrees of arthritis of the control animals and of the compound-treated animals.

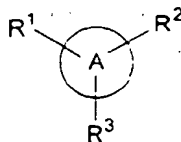
[0714] From the rates of treatment and the dosages, ID₅₀ values were calculated by the least squares method.

[0715] In this test, the compounds of the present invention showed excellent activity in treating arthritis induced by anti-collagen antibody.

[0716] As illustrated above, the compounds of the present invention exhibit excellent activity in inhibiting the production of inflammatory cytokines, particularly in inhibiting the production of IL-1 β and TNF α . Furthermore, the compounds of the present invention have satisfactory oral absorptivity and a low level of toxicity. Consequently, the compounds of the present invention are useful as pharmaceuticals, suitable for the prophylaxis and treatment of both humans and animals. They can, for example, be used as an analgesic, an anti-inflammatory agent and an antiviral agent as well as an agent for use in the prophylaxis and treatment of chronic rheumatoid arthritis, degenerative arthritis, allergic diseases, asthma, septicaemia, psoriasis, osteoporosis, autoimmune diseases (e.g., systemic lupus erythematosus, ulcerative colitis and Crohn's disease), diabetes, glomerular nephritis, hepatitis and arteriosclerosis. Of these applications, the compounds of the present invention are particularly useful as an analgesic and an anti-inflammatory agent and as an agent for the prophylaxis and treatment of chronic rheumatoid arthritis, degenerative arthritis, allergic diseases, septicaemia, psoriasis, osteoporosis, ulcerative colitis, diabetes, hepatitis and arteriosclerosis.

Claims

1. A compound of formula (I), or a pharmacologically acceptable salt, ester or other derivative thereof:



(I)

wherein:

A represents a pyrrole ring;

R¹ is an aryl group defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below, or a heteroaryl group defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below;

R² represents a heteroaryl group defined below having at least one ring nitrogen atom, said heteroaryl group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below, Substituent group β defined below, Substituent group γ defined below and Substituent group δ defined below;

R³ represents a group of the formula -X-R⁴, wherein:

X is selected from the group consisting of single bonds, lower alkylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below, lower alkenylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below, and lower alkynylene groups, defined below, which may optionally be substituted with at least one substituent selected from Substituent group α defined below; and

R⁴ is selected from the group consisting of

cycloalkyl groups defined below which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below, aryl groups defined below which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below, heterocyclyl groups defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below, heterocyclyl groups defined below having at least one nitrogen atom, said heterocyclyl groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below, heteroaryl groups defined below which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below

and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined below and Substituent group γ defined below, heteroaryl groups defined below having at least one nitrogen atom, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α defined below and Substituent group δ defined below, and groups of formula $-NR^aR^b$, wherein R^a and R^b are the same or different from each other and each is independently selected from the group consisting of hydrogen atoms, lower alkyl groups defined below, lower alkenyl groups defined below, lower alkynyl groups defined below, aralkyl groups defined below and lower alkylsulfonyl groups defined below;

PROVIDED THAT said substituents R^1 and R^3 are bonded to the two atoms of said pyrrole ring which are adjacent to the atom of the pyrrole ring to which said substituent R^2 is bonded;

Substituent group α comprises hydroxyl groups, nitro groups, cyano groups, halogen atoms, lower alkoxy groups defined below, halogeno lower alkoxy groups defined below, lower alkylthio groups defined below, and halogeno lower alkylthio groups defined below;

Substituent group β comprises groups of formula $-NR^cR^d$, wherein R^c and R^d are the same or different from each other and each is independently selected from the group consisting of hydrogen atoms, lower alkyl groups defined below, lower alkenyl groups defined below, lower alkynyl groups defined below, aralkyl groups defined below and lower alkylsulfonyl groups, or R^c and R^d together with the nitrogen atom to which R^c and R^d are bonded, form a heterocyclyl group defined below;

Substituent group γ comprises lower alkyl groups defined below which are substituted with a group of formula $-NR^cR^d$ wherein R^c and R^d are as defined above; and Substituent group δ comprises lower alkyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, lower alkenyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, lower alkynyl groups defined below which may optionally be substituted with at least one substituent selected from Substituent group α defined above, aralkyl groups defined below and cycloalkyl groups defined below;

said optionally substituted aryl groups in the definitions of substituents R^1 and R^4 above are aromatic hydrocarbon groups having from 6 to 14 carbon atoms in one or more rings, said aromatic hydrocarbon groups optionally being fused with a cycloalkyl group having from 3 to 10 carbon atoms;

said optionally substituted heteroaryl groups in the definition of substituent R^1 above and said heteroaryl groups which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above and which may optionally be further substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above in the definition of substituent R^4 above are 5-to 7-membered aromatic heterocyclic groups containing from 1 to 3 heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, said aromatic heterocyclic groups optionally being fused with another cyclic group selected from the group consisting of aryl groups as defined above and cycloalkyl groups having from 3 to 10 carbon atoms;

said optionally substituted heteroaryl groups having at least one ring nitrogen atom in the definition of substituent R^2 above and said optionally substituted heteroaryl groups having at least one ring nitrogen atom in the definition of substituent R^4 above are 5- to 7-membered aromatic heterocyclic groups containing one or two nitrogen atoms and optionally containing one or two further heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms;

said lower alkylene groups in the definition of group X above are straight or branched alkylene groups having from 1 to 6 carbon atoms;

said lower alkenylene groups in the definition of group X above are straight or branched alkenylene groups having from 2 to 6 carbon atoms;

said lower alkynylene groups in the definition of group X above are straight or branched alkynylene groups having from 2 to 6 carbon atoms;

said cycloalkyl groups in the definitions of substituent R⁴ above and Substituent group δ above are cycloalkyl groups having from 3 to 7 carbon atoms;

said heterocyclyl groups which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above and which are substituted with at least one substituent selected from the group consisting of Substituent group β defined above and Substituent group γ defined above in the definition of substituent R⁴ above are non-aromatic heterocyclic groups having from 4 to 14 ring atoms in one or more rings, at least one of said ring atoms being a heteroatom selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, said heterocyclyl groups optionally being fused with another cyclic group selected from the group consisting of aryl groups as defined above and heteroaryl groups as defined above;

said heterocyclyl groups having at least one ring nitrogen atom in which said heterocyclyl groups may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α defined above and Substituent group δ defined above in the definition of substituent R⁴ above are 4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally containing one further heteroatom selected from the group consisting of oxygen atoms, sulfur atoms and nitrogen atoms, said non-aromatic heterocyclic groups optionally being fused with another cyclic group selected from the group consisting of aryl groups as defined above and heteroaryl groups as defined above;

said lower alkyl groups in the definition of substituents R^a, R^b, R^c and R^d above, said lower alkyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ above and said lower alkyl moiety of the lower alkyl groups substituted with a group of formula -NR^cR^d in the definition of Substituent group γ above are straight or branched alkyl groups having from 1 to 6 carbon atoms;

said lower alkenyl groups in the definition of substituents R^a, R^b, R^c and R^d above and said lower alkenyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ above are straight or branched alkenyl groups having from 2 to 6 carbon atoms;

said lower alkynyl groups in the definition of substituents R^a, R^b, R^c and R^d above and said lower alkynyl groups which may optionally be substituted with at least one substituent selected from Substituent group α in the definition of Substituent group δ above are straight or branched alkynyl groups having from 2 to 6 carbon atoms;

said aralkyl groups in the definition of R^a, R^b, R^c and R^d above and in the definition of Substituent group δ above are lower alkyl groups as defined above which are substituted with at least one aryl group as defined above;

said lower alkylsulfonyl groups in the definition of substituents R^a, R^b, R^c and R^d are lower alkyl groups as defined above which are bonded to a sulfonyl group;

said heterocyclyl groups formed by substituents R^c and R^d together with the nitrogen atom to which they are bonded are heterocyclyl groups having at least one nitrogen atom as defined above in relation to substituent R⁴ above, said heterocyclyl groups optionally being fused with another cyclic group selected from the group consisting of aryl groups as defined above and heteroaryl groups as defined above;

said lower alkoxy groups in the definition of Substituent group α above are lower alkyl groups as defined above which are bonded to an oxygen atom;

said halogeno lower alkoxy groups in the definition of Substituent group α above are lower alkoxy groups as defined above which are substituted with at least one halogen atom;

said lower alkylthio groups in the definition of Substituent group α above are lower alkyl groups as defined

above which are bonded to a sulfur atom; and

said halogeno lower alkylthio groups in the definition of Substituent group α above are lower alkylthio groups as defined above which are substituted with at least one halogen atom.

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2. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^1 is an aryl group as defined in claim 1 which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

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3. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^1 is a phenyl or naphthyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

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4. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^1 is a phenyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α^1 defined below, Substituent group β^1 defined below, Substituent group γ^1 defined below and Substituent group δ^1 defined below;

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said Substituent group α^1 comprises halogen atoms, lower alkoxy groups as defined in claim 1 and halogeno lower alkoxy groups as defined in claim 1;

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said Substituent group β^1 comprises a group of formula $-NR^cR^d$ wherein one of R^c and R^d is selected from the group consisting of hydrogen atoms and lower alkyl groups as defined in claim 1 and the other is selected from the group consisting of hydrogen atoms, lower alkyl groups as defined in claim 1 and aralkyl groups as defined in claim 1;

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said Substituent group γ^1 comprises lower alkyl groups as defined in claim 1 which are substituted with a substituent selected from the group consisting of amino groups, amino groups substituted with one or two lower alkyl groups as defined in claim 1 and amino groups substituted with an aralkyl group as defined in claim 1; and

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said Substituent group δ^1 comprises lower alkyl groups as defined in claim 1, halogeno lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one halogen atom, hydroxy lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one hydroxy group and nitro lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one nitro group.

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5. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^1 is a phenyl group which may optionally be substituted with at least one substituent selected from the group consisting of halogen atoms, halogeno lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one halogen atom and halogeno lower alkoxy groups as defined in claim 1.

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6. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^1 is selected from the group consisting of phenyl, 4-fluorophenyl, 3-fluorophenyl, 3-chlorophenyl, 3,4-difluorophenyl, 3,4,5-trifluorophenyl, 3-chloro-4-fluorophenyl, 3-difluoromethoxyphenyl and 3-trifluoromethylphenyl groups.

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7. A compound according to any one of claims 1 to 6 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^2 is a 5- or 6-membered aromatic heterocyclic group containing one or two nitrogen atoms, said group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

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8. A compound according to any one of claims 1 to 6 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^2 is a pyridyl or pyrimidinyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

9. A compound according to any one of claims 1 to 6 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted with at least one

substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

10. A compound according to any one of claims 1 to 6 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

11. A compound according to any one of claims 1 to 6 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of methoxy, amino, methylamino, benzylamino and α -methylbenzylamino groups.

12. A compound according to any one of claims 1 to 11 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^4 is selected from the group consisting of:

cycloalkyl groups having from 3 to 7 carbon atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

phenyl or naphthyl groups which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally one further heteroatom selected from the group consisting of oxygen atoms, sulfur atom and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ ;

5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

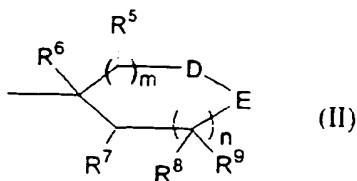
5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ ; and groups of formula $-NR^aR^b$ wherein R^a and R^b are as defined in claim 1.

13. A compound according to any one of claims 1 to 12 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein X is selected from the group consisting of single bonds, alkylene groups having from 1 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α , alkenylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α and alkynylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α .

14. A compound according to any one of claims 1 to 12 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein X is selected from the group consisting of single bonds, alkylene groups having from 1 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α , alkenylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α and alkynylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α .

15. A compound according to any one of claims 1 to 12 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein X is a single bond or an alkenylene group having from 2 to 4 carbon atoms.

16. A compound according to any one of claims 1 to 11 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^3 is a group of the following formula (II):



wherein:

m represents 0 or 1;

n represents 1 or 2;

R^5 , R^6 , R^7 , R^8 and R^9 may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ;

one of D and E represents a group of formula $>NR^{10}$ wherein R^{10} is selected from the group consisting of hydrogen atoms, Substituent group γ and Substituent group δ , and the other represents a group of formula $>CR^{11}R^{12}$ wherein R^{11} and R^{12} are the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , or

R^6 may form, together with R^5 or R^7 , a single bond, and/or R^{10} and R^{11} together may form a straight or branched alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

17. A compound according to claim 16 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein m is 1 and n is 1.

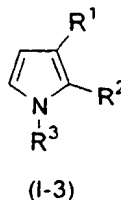
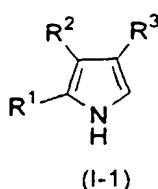
18. A compound according to claim 16 or claim 17 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein E represents a group of formula $>NR^{10}$ and D represents a group of formula $>CR^{11}R^{12}$.

19. A compound according to any one of claims 16 to 18 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein one of R^5 and R^7 together with R^6 represents a single bond and the other represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , R^8 , R^9 and R^{12} may be the same or different from one another and each represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , and R^{10} and R^{11} may be the same or different from one another and each represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ or R^{10} and R^{11} together form an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

20. A compound according to any one of claims 16 to 19 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein R^{10} and R^{11} together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

21. A compound according to any one of claims 16 to 20 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein each of R^7 , R^8 and R^9 is a hydrogen atom, and R^{10} and R^{11} together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

22. A compound according to any one of claims 1 to 21 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein said compound of formula (I) is a compound of formula (I-1) or (I-3) below:



23. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R¹ is an aryl group as defined in claim 1 which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ;

R² is a 5- or 6-membered aromatic heterocyclic group containing one or two nitrogen atoms, said group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ; and

R³ is a group of formula -X-R⁴ wherein:

R⁴ is selected from the group consisting of

cycloalkyl groups having from 3 to 7 carbon atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

phenyl or naphthyl groups which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally one further heteroatom selected from the group consisting of oxygen atoms, sulfur atom and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ ;

5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ , and groups of formula -NR^aR^b wherein R^a and R^b are as defined in claim 1; and

X is selected from the group consisting of single bonds, alkylene groups having from 1 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α , alkynylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α and alkynylene groups having from 2 to 6 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α .

24. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R¹ is an aryl group as defined in claim 1 which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Sub-

stituent group δ ;

R^2 is a 5- or 6-membered aromatic heterocyclic group containing one or two nitrogen atoms, said group optionally being substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ; and

R^3 is a group of formula $-X-R^4$ wherein:

R^4 is selected from the group consisting of

cycloalkyl groups having from 3 to 7 carbon atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

phenyl or naphthyl groups which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally one further heteroatom selected from the group consisting of oxygen atoms, sulfur atom and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ ;

5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ ;

5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ , and groups of formula $-NR^aR^b$ wherein R^a and R^b are as defined in claim 1; and

X is selected from the group consisting of single bonds, alkylene groups having from 1 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α , alkenylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α and alkynylene groups having from 2 to 4 carbon atoms which may optionally be substituted with at least one substituent selected from Substituent group α .

25. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R^1 is a phenyl or naphthyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ;

R^2 is a pyridyl or pyrimidinyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ; and

R^3 is a group of formula $-X-R^4$ wherein

R^4 is selected from the group consisting of

cycloalkyl groups having from 3 to 7 carbon atoms, which may optionally be substituted with at least

one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ .

phenyl or naphthyl groups which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ .

4- to 12-membered non-aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ .

4- to 12-membered non-aromatic heterocyclic groups containing one nitrogen atom and optionally one further heteroatom selected from the group consisting of oxygen atoms, sulfur atom and nitrogen atoms, which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ .

5- or 6-membered aromatic heterocyclic groups containing one or two heteroatoms selected from the group consisting of sulfur atoms, oxygen atoms and nitrogen atoms, which may be optionally substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ and which are substituted with at least one substituent selected from the group consisting of Substituent group β and Substituent group γ .

5- or 6-membered aromatic heterocyclic groups containing one or two nitrogen atoms, said groups optionally being substituted with at least one substituent selected from the group consisting of Substituent group α and Substituent group δ , and groups of formula $-NR^aR^b$ wherein R^a and R^b are as defined in claim 1; and

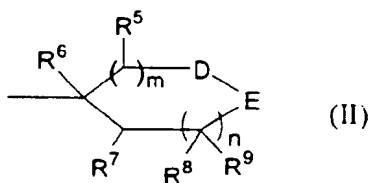
X is a single bond or an alkenylene group having from 2 to 4 carbon atoms.

26. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R^1 is a phenyl or naphthyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ;

R^2 is a pyridyl or pyrimidinyl group which may optionally be substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ; and

R^3 is a group of the following formula (II):



wherein

m represents 0 or 1,

n represents 1 or 2,

R^5 , R^6 , R^7 , R^8 and R^9 may be the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group β , Substituent group γ and Substituent group δ , one of D and E represents a group of formula $>NR^{10}$ wherein R^{10} is selected from the group consisting of hydrogen atoms, Substituent group γ and Substituent group δ , and the other represents a group of formula $>CR^{11}R^{12}$ wherein R^{11} and R^{12} are the same or different from one another and each is selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

uent group γ and Substituent group δ , or
 R^6 may form, together with R^5 or R^7 , a single bond, and/or R^{10} and R^{11} together may form a straight or
 branched alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at
 least one substituent selected from the group consisting of Substituent group α , Substituent group β ,
 Substituent group γ and Substituent group δ .

27. A compound according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R^1 is a phenyl group which may optionally be substituted with at least one substituent selected from the group
 consisting of Substituent group α^1 defined below, Substituent group β^1 defined below, Substituent group γ^1
 defined below and Substituent group δ^1 defined below;

R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted with at least one substituent se-
 lected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Sub-
 stituent group δ ; and

R^3 is a group of formula (II) as defined in claim 16 wherein m is 1 and n is 1;

said Substituent group α^1 comprises halogen atoms, lower alkoxy groups as defined in claim 1 and halogeno
 lower alkoxy groups as defined in claim 1;

said Substituent group β^1 comprises a group of formula $-NR^cR^d$ wherein one of R^c and R^d is selected from
 the group consisting of hydrogen atoms and lower alkyl groups as defined in claim 1 and the other is selected
 from the group consisting of hydrogen atoms, lower alkyl groups as defined in claim 1 and aralkyl groups as
 defined in claim 1;

said Substituent group γ^1 comprises lower alkyl groups as defined in claim 1 which are substituted with a
 substituent selected from the group consisting of amino groups, amino groups substituted with one or two
 lower alkyl groups as defined in claim 1 and amino groups substituted with an aralkyl group as defined in claim
 1; and

said Substituent group δ^1 comprises lower alkyl groups as defined in claim 1, halogeno lower alkyl groups
 comprising lower alkyl groups as defined in claim 1 which are substituted with at least one halogen atom,
 hydroxy lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at
 least one hydroxy group and nitro lower alkyl groups comprising lower alkyl groups as defined in claim 1 which
 are substituted with at least one nitro group.

28. A compound of formula (I) according to claim 1 or a pharmacologically acceptable salt, ester or other derivative
 thereof, wherein:

R^1 is a phenyl group which may optionally be substituted with at least one substituent selected from the group
 consisting of Substituent group α^1 defined below, Substituent group β^1 defined below, Substituent group γ^1
 defined below and Substituent group δ^1 defined below;

R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted with at least one substituent se-
 lected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substit-
 uent group δ ; and

R^3 is a group of formula (II) as defined in claim 16 wherein E represents a group of formula $>NR^{10}$ wherein
 R^{10} is as defined in claim 1 and D represents a group of formula $>CR^{11}R^{12}$ wherein R^{11} and R^{12} are as defined
 in claim 1;

said Substituent group α^1 comprises halogen atoms, lower alkoxy groups as defined in claim 1 and halogeno
 lower alkoxy groups as defined in claim 1;

said Substituent group β^1 comprises a group of formula $-NR^cR^d$ wherein one of R^c and R^d is selected from
 the group consisting of hydrogen atoms and lower alkyl groups as defined in claim 1 and the other is selected
 from the group consisting of hydrogen atoms, lower alkyl groups as defined in claim 1 and aralkyl groups as

defined in claim 1;

said Substituent group γ^1 comprises lower alkyl groups as defined in claim 1 which are substituted with a substituent selected from the group consisting of amino groups, amino groups substituted with one or two lower alkyl groups as defined in claim 1 and amino groups substituted with an aralkyl group as defined in claim 1; and

said Substituent group δ^1 comprises lower alkyl groups as defined in claim 1, halogeno lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one halogen atom, hydroxy lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one hydroxy group and nitro lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one nitro group.

29. A compound of formula (I) according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein

R^1 is a phenyl group which may optionally be substituted with at least one substituent selected from the group consisting of halogen atoms, halogeno lower alkyl groups comprising lower alkyl groups as defined in claim 1 which are substituted with at least one halogen atom and halogeno lower alkoxy groups as defined in claim 1;

R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ ; and

R^3 is a group of formula (II) as defined in claim 16 wherein one of R^5 and R^7 together with R^6 represents a single bond and the other represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , R^8 , R^9 and R^{12} may be the same or different from one another and each represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ , and R^{10} and R^{11} may be the same or different from one another and each represents a substituent selected from the group consisting of hydrogen atoms, Substituent group α , Substituent group γ and Substituent group δ or R^{10} and R^{11} together form an alkylene group having from 1 to 6 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

30. A compound of formula (I) according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R^1 is selected from the group consisting of phenyl, 4-fluorophenyl, 3-fluorophenyl, 3-chlorophenyl, 3,4-difluorophenyl, 3,4,5-trifluorophenyl, 3-chloro-4-fluorophenyl, 3-difluoromethoxyphenyl and 3-trifluoromethylphenyl groups;

R^2 is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of methoxy, amino, methylamino, benzylamino and α -methylbenzylamino groups; and

R^3 is a group of formula (II) as defined in claim 16 wherein R^{10} and R^{11} together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

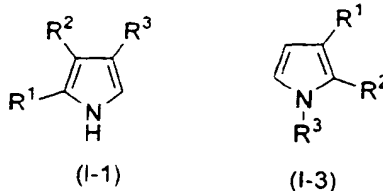
31. A compound of formula (I) according to claim 1 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein:

R^1 is selected from the group consisting of phenyl, 4-fluorophenyl, 3-fluorophenyl, 3-chlorophenyl, 3,4-difluorophenyl, 3,4,5-trifluorophenyl, 3-chloro-4-fluorophenyl, 3-difluoromethoxyphenyl and 3-trifluoromethylphenyl groups;

R² is a 4-pyridyl or 4-pyrimidinyl group which may optionally be substituted at the 2-position thereof with a substituent selected from the group consisting of methoxy, amino, methylamino, benzylamino and α -methylbenzylamino groups; and

R³ is a group of formula (II) as defined in claim 16 wherein each of R⁷, R⁸ and R⁹ is a hydrogen atom, and R¹⁰ and R¹¹ together form a straight chain alkylene group having 3 or 4 carbon atoms which is unsubstituted or is substituted with at least one substituent selected from the group consisting of Substituent group α , Substituent group β , Substituent group γ and Substituent group δ .

32. A compound of formula (I) according to any one of claims 23 to 31 or a pharmacologically acceptable salt, ester or other derivative thereof, wherein said compound of formula (I) is a compound of formula (I-1) or (I-3) below:



33. A compound of formula (I) according to claim 1 selected from the following group of compounds, or a pharmacologically acceptable salt, ester or other derivative thereof:

4-(3-aminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(3-acetylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(3-methylaminopropyl)-3-(pyridin-4-yl)-1H-pyrrole,
 [5-(4-fluorophenyl)-4-(pyridin-4-yl)-1H-pyrrol-3-yl]-(pyridin-4-yl)methanol,
 4-(1-acetyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(piperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1H-pyrrole,
 1-(1-acetyl-piperidin-4-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-1-[1-(2-nitroethyl)piperidin-4-yl]-2-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-3-yl)-2-(pyridin-4-yl)-1H-pyrrole,
 1-(azetidin-3-yl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole,
 4-(3-dimethylaminopropyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-[3-(piperidin-1-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-[3-(1-methylpiperazin-4-yl)propyl]-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(1-methylpiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(2,2,6,6-tetramethylpiperidin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-2-en-3-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(quinuclidin-3-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(4-hydroxypiperidin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-2-(pyridin-4-yl)-1-(quinuclidin-3-yl)-1H-pyrrole,
 1-(4-aminocyclohexyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1H-pyrrole,
 3-(4-fluorophenyl)-2-(2-methylaminopyrimidin-4-yl)-1-(piperidin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]oct-2-en-3-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(8-methyl-8-azabicyclo[3.2.1]octan-3-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydro-2,2,6,6-tetramethylpyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(2,2,6,6-tetramethylpiperidin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1H-pyrrole,
 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1H-pyrrole,

4-(1-ethylpiperidin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,6-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,2,6,6-pentamethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(3-dimethylamino-1-propen-1-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(4-aminobutyl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 1-(3-dimethylaminopropyl)-3-(4-fluorophenyl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,7,8,8a-octahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(6-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(2-allyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(6-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(2-benzyl-1-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,9,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(2-methylaminopyrimidin-4-yl)-1 *H*-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1 *H*-pyrrole,
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1 *H*-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3-chlorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2,3,5,8,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1 *H*-pyrrole, and
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-2-phenyl-3-(pyridin-4-yl)-1 *H*-pyrrole.

34. A compound of formula (I) according to claim 1 selected from the following group of compounds, or a pharmacologically acceptable salt, ester or other derivative thereof:

2-(4-fluorophenyl)-3-(pyridin-4-yl)-4-(1,2,3,6-tetrahydropyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-(piperidin-4-yl)-2-(pyridin-4-yl)-1 *H*-pyrrole,
 3-(4-fluorophenyl)-1-[3-(morpholin-1-yl)propyl]-2-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,6-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1,2-dimethyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-6-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 4-(1-ethyl-2-methyl-1,2,3,6-tetrahydropyridin-4-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2*H*-quinolizin-8-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole,
 2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1 *H*-pyrrole, and
 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1 *H*-pyrrole,

35. A compound according to claim 1 selected from 4-(8-azabicyclo[3.2.1]oct-2-en-3-yl)-2-(4-fluorophenyl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
36. A compound according to claim 1 selected from 2-(4-fluorophenyl)-4-(1-phenethyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
37. A compound according to claim 1 selected from 2-(4-fluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
38. A compound according to claim 1 selected from 2-(4-fluorophenyl)-4-(6-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
39. A compound according to claim 1 selected from 2-(4-fluorophenyl)-4-(2-methyl-1-propyl-1,2,3,6-tetrahydropyridin-4-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
40. A compound according to claim 1 selected from 2-(4-fluorophenyl)-4-(1,3,4,6,7,9a-hexahydro-2H-quinolizin-8-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
41. A compound according to claim 1 selected from 2-(3,4-difluorophenyl)-4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
42. A compound according to claim 1 selected from 4-(1,2,3,5,6,8a-hexahydroindolizin-7-yl)-3-(pyridin-4-yl)-2-(3-trifluoromethylphenyl)-1*H*-pyrrole, and pharmacologically acceptable salts, esters and other derivatives thereof.
43. A pharmaceutical composition comprising an effective amount of a pharmacologically active compound together with a carrier or diluent therefor, wherein said pharmacologically active compound is a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof.
44. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use as a medicament.
45. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in inhibiting the production of inflammatory cytokines in a mammal, which may be human.
46. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in inhibiting bone resorption.
47. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of inflammatory diseases.
48. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of viral diseases.
49. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in relieving pain or pyrexia.
50. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of chronic rheumatoid arthritis.
51. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of osteoarthritis.
52. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of cancer.
53. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of hepatitis.

54. A compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof for use in the treatment or prophylaxis of a disease selected from the group consisting of allergic diseases, septicæmia, psoriasis, osteoporosis, ulcerative colitis, diabetes, nephritis, ischemic heart disease, Alzheimer's disease and arteriosclerosis.
55. The use of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of diseases which can be treated or prevented by inhibiting the production of inflammatory cytokines in a mammal, which may be human.
56. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for inhibiting bone resorption.
57. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of inflammatory diseases.
58. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of viral diseases.
59. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for relieving pain or pyrexia.
60. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of chronic rheumatoid arthritis.
61. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of osteoarthritis.
62. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of cancer.
63. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of hepatitis.
64. The use according to claim 55 of a compound of formula (I) according to any one of claims 1 to 42 or a pharmacologically acceptable salt, ester or other derivative thereof in the manufacture of a medicament for the treatment or prophylaxis of a disease selected from the group consisting of allergic diseases, septicæmia, psoriasis, osteoporosis, ulcerative colitis, diabetes, nephritis, ischemic heart disease, Alzheimer's disease and arteriosclerosis.

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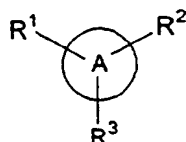
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(54) **Heteroaryl-substituted pyrrole derivatives, their preparation and their therapeutic uses**

(57) Compounds of formula (I):



(I)

[wherein: A is a pyrrole ring; R¹ is an optionally substituted aryl or heteroaryl group; R² is an optionally substituted nitrogen-containing heteroaryl group; R³ repre-

sents a group of the formula -X-R⁴, wherein X is a single bond or an optionally substituted alkylene, alkenylene or alkynylene group, and R⁴ is a substituted cycloalkyl group, a substituted aryl group, a substituted heterocyclyl group, an optionally substituted nitrogen-containing heterocyclyl group, a substituted heteroaryl group, an optionally substituted nitrogen-containing heteroaryl group, or -NR^aR^b, wherein each of R^a and R^b is a hydrogen atom or an alkyl, alkenyl, alkynyl, aralkyl or alkylsulfonyl group; PROVIDED THAT said substituents R¹ and R³ are bonded to the two atoms of said pyrrole ring which are adjacent to the atom of the pyrrole ring to which said substituent R² is bonded] have excellent inhibitory activity against the production of inflammatory cytokines.

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Office

PARTIAL EUROPEAN SEARCH REPORT

Application Number

which under Rule 45 of the European Patent Convention shall be considered, for the purposes of subsequent proceedings, as the European search report

EP 00 30 6196

DOCUMENTS CONSIDERED TO BE RELEVANT			
Category	Citation of document with indication, where appropriate, of relevant passages	Relevant to claim	CLASSIFICATION OF THE APPLICATION (Int.Cl.7)
Y,D	WO 97 05877 A (MERCK & CO INC ; LASZLO STEPHEN E DE (US); MANTLO NATHAN B (US); P0) 20 February 1997 (1997-02-20) * claims 1,26-43; example 4 *	1,43-64	C07D401/04 C07D401/14 C07D453/02 C07D451/02 C07D471/04 C07D455/02
Y,D	WO 96 21452 A (ADAMS JERRY LEROY ; BOEHM JEFFREY CHARLES (US); GALLAGHER TIMOTHY F) 18 July 1996 (1996-07-18) * claims; example 23 *	1,43-64	A61P11/06 A61P9/10 A61P17/06 A61P29/00
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INCOMPLETE SEARCH			
<p>The Search Division considers that the present application, or one or more of its claims, does/do not comply with the EPC to such an extent that a meaningful search into the state of the art cannot be carried out, or can only be carried out partially, for these claims.</p> <p>Claims searched completely:</p> <p>Claims searched incompletely:</p> <p>Claims not searched:</p> <p>Reason for the limitation of the search</p> <p>see sheet C</p>			
Place of search		Date of completion of the search	Examiner
THE HAGUE		27 November 2000	Bosma, P
CATEGORY OF CITED DOCUMENTS			
<p>X: particularly relevant if taken alone</p> <p>Y: particularly relevant if combined with another document of the same category</p> <p>A: technological background</p> <p>O: non-written disclosure</p> <p>P: intermediate document</p>		<p>T: theory or principle underlying the invention</p> <p>E: earlier patent document, but published on, or after the filing date</p> <p>D: document cited in the application</p> <p>L: document cited for other reasons</p> <p>&: member of the same patent family, corresponding document</p>	

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European Patent
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INCOMPLETE SEARCH
SHEET C

Application Number
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Claim(s) searched completely:
35-42

Claim(s) searched incompletely:
1-34, 43-64

Reason for the limitation of the search:

Present claims 1-34, and 43-64 relate to an extremely large number of possible compounds. Support within the meaning of Article 84 EPC and/or disclosure within the meaning of Article 83 EPC is to be found, however, for only a very small proportion of the compounds claimed. In the present case, the claims so lack support, and the application so lacks disclosure, that a meaningful search over the whole of the claimed scope is impossible. Consequently, the search has been carried out for those parts of the claims which appear to be supported and disclosed, namely those parts relating to the compounds of formula (I) in which R1 has the meanings defined by claim 6, and R2 has the meanings defined by claim 8.

**ANNEX TO THE EUROPEAN SEARCH REPORT
ON EUROPEAN PATENT APPLICATION NO.**

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This annex lists the patent family members relating to the patent documents cited in the above-mentioned European search report. The members are as contained in the European Patent Office EDP file on
The European Patent Office is in no way liable for these particulars which are merely given for the purpose of information.

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For more details about this annex : see Official Journal of the European Patent Office, No. 12/82